

PILOT STUDY WORK PLAN

West Coast Door Site Tacoma, Washington

January 13, 2025

Prepared for

3102 Tenants in Common 3102/3120 South Pine Street Tacoma, Washington

Pilot Study Work Plan West Coast Door Site Tacoma, Washington

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LIST OF ABBREVIATIONS AND ACRONYMS

AO	Agreed Order
	aboveground storage tank
bgs	below ground surface
	benzene, toluene, ethylbenzene, xylenes
CFM	cubic feet per minute
cPAH	carcinogenic PAH
CSL	Contaminated Sites List
CSM	conceptual site model
CUL	cleanup level
DNAPL	dense nonaqueous phase liquid
DO	dissolved oxygen
Ecology	Washington State Department of Ecology
EISB	enhanced in situ bioremediation
EPA	US Environmental Protection Agency
FS	feasibility study
ft	foot/feet
gpm	gallons per minute
HASP	Health and Safety Plan
IDW	investigation-derived waste
IET	Innovative Environmental Technologies, Inc.
ISCO	in situ chemical oxidation
ISGS	in situ geochemical stabilization
	Landau Associates, Inc.
μg/m ³	micrograms per cubic meter
μg/L	micrograms per liter
mg/kg	milligrams per kilogram
MTCA	Model Toxics Control Act
MW	monitoring well
•	North American Datum of 1983/1991
NAPL	non-aqueous phase liquid
NAVD88	North American Vertical Datum of 1988
PAH	polycyclic aromatic hydrocarbon
	Pacific Groundwater Group
PPE	personal protective equipment
·	pounds per square inch
psig	pounds per square inch gauge
	polyvinyl chloride
RCRA	Resource Conservation and Recovery Act

LIST OF ABBREVIATIONS AND ACRONYMS (continued)

RI	remedial investigation
ROI	radius of influence
ROW	right-of-way
SF	square feet
SIM	selected ion monitoring
Site	West Coast Door Site
SVOC	semivolatile organic compound
TCE	trichloroethene
TEQ	toxic equivalent
TIC	Tenants in Common
TPH	total petroleum hydrocarbon
UIC	underground injection control
UST	underground storage tank
VOC	volatile organic compound
WAC	Washington Administrative Code
work plan	pilot study work plan

1.0 INTRODUCTION

Landau Associates, Inc. (Landau) has prepared this pilot study work plan (work plan) to assist the West Coast Door Site (Site) former owners (collectively referred to as 3102 Tenants in Common [TIC]) with the remediation of the Site located at 3133 South Cedar Street in Tacoma, Washington. The work plan has been prepared in accordance with an Agreed Order (AO; No. DE 14016, April 11, 2017) with the Washington State Department of Ecology (Ecology) that requires 3102 TIC to address contamination resulting from historical creosote-treated wood pipe manufacturing operations. The Site is included on Ecology's Contaminated Sites List (CSL) under Facility Site ID 6308485 and Cleanup Site ID 2599.

1.1 Site Description

The Site property consists of two irregularly shaped parcels (Pierce County parcel numbers 0320073069 and 0320073070) covering 10.43 acres of land in the City of Tacoma, Pierce County, Washington (Figure 1). It is bounded to the north by South Center Street, to the west by South Cedar Street, to the south by Sound Transit-owned railroad tracks, and to the east by South Pine Street (Figure 2). The property is zoned for industrial use, and its western portion is occupied by adjoining north and south warehouse buildings constructed in the mid-1980s that cover approximately 89,000 and 108,000 square feet (SF), respectively. The eastern portion consists of loading docks adjacent to the warehouse buildings and a paved parking lot. The warehouse buildings on the western portion of the Site are leased to Goodwill Industries, which uses the north warehouse for shipping, receiving, and sorting operations. A section of the south warehouse is used as a retail outlet that is open to the public.

Approximately 95 percent of the surface area of the property is capped with impermeable surfaces, including pavement and structures. The remaining 5 percent of uncapped surfaces consists of grass areas to the west of the parking area on the southern end of the southern warehouse buildings. The rights-of-way (ROWs) surrounding the property to the north, west, and east are also paved; however, the Sound Transit-owned ROW to the south of the Site is partially unpaved and surfaced with compacted crushed gravel. Surrounding property uses are primarily warehousing or industrial operations.

1.2 Site Background

The Site was originally developed by Buffelen Pipe and Creosote Company and American Wood Pipe Company, which operated from the early 1900s to the mid-1930s. Manufacturing operations included log storage, drying kilns, and a creosoting plant that included two pressurized retort vessels used to impregnate pipe with creosote and was located in the southwestern portion of the current south warehouse footprint. The creosote oil consisted of a distillate of coal-gas tar or coke-oven tar and was stored outside in creosote oil tanks that had a combined capacity of 50,000 gallons.

Monarch Door and Manufacturing Company began door manufacturing at the Site in the mid-1930s. West Coast Door, Inc. purchased the subject property in 1954. West Coast Door manufacturing operations included cutting, sanding, and gluing of wood-veneered fiberboard core doors. These operations continued in both the north and south portions of the current warehouse after they were

constructed in the mid-1980s. William B. Swensen (also operating as William B. Swensen Enterprises) purchased the property in the 1970s and operated it as West Coast Door until 1997. Door manufacturing operations ceased in 1997, at which point the facilities were converted for warehouse and office space use. Tenants in the southern warehouse included Total Recall Information Management, Thrifty Supply, Nalley Fine Foods, and Goodwill Industries. Ownership of the Site was passed to the heirs of William B. Swensen, operating as 3102 TIC, after Mr. Swensen's death in 2006. CenterPoint Properties Trust subsequently purchased the property in 2017 and is the current owner.

In addition to the current warehouse structure, five additional smaller buildings were constructed between 1961 and 1970 on the central and eastern portion of the subject property and were demolished after 2006. These structures included a warehouse/showroom, office buildings, a maintenance shop, and a shipping/wood storage building.

Excavation and grading activities undertaken in 1986 as part of construction of the south warehouse revealed the presence of creosote-like material in subsurface soils. Approximately 10,500 pounds of material containing greater than 1 percent polycyclic aromatic hydrocarbons (PAHs) and designated as Washington State extremely hazardous waste were excavated and disposed offsite under a Resource Conservation and Recovery Act (RCRA) dangerous waste permit. The north warehouse, completed in 1985, and south warehouse, completed in 1987, are both occupied by Goodwill Industries.

Three storage tanks containing gasoline have previously existed on the Site, including two underground storage tanks (USTs) of 3,000- and 2,500-gallon capacity, which were removed from the Site in 1989. Confirmation soil samples from this tank removal were submitted to the Tacoma-Pierce County Health department, which certified the tank removal as complete based on confirmation results less than required cleanup levels (CULs). A third tank, an aboveground storage tank (AST) installed in 1990, was removed from the Site in 2005, and soil sampling conducted in the vicinity confirmed that gasoline was not released to soil.

1.3 Natural Conditions

Natural conditions at the Site, including the physical setting, geology, and hydrogeology, are discussed in the subsections below.

1.3.1 Physical Setting

The Site is situated in the South Tacoma Channel of the Nalley Valley in Tacoma. The Site is generally flat, with an elevation of approximately 250 feet (ft) NAVD88.¹

1.3.2 Geology

The Site lies in the South Tacoma Channel, which is filled by Vashon-era fluvial deposits that were deposited in a high-energy glaciofluvial environment and are composed primarily of sand and gravel. The Vashon-era deposits were created as a result of the South Tacoma channel acting as a spillway for proglacial lakes that formed during the recession of the Vashon ice sheet in the late Pleistocene. The

¹ North American Vertical Datum of 1988.

valley is one of the major channels that connected glacial Lake Puyallup through progressively lower spillways into Lake Russell, the main proglacial lake in front of the receding Vashon Ice sheet (Troost and Sofield 2011; Walters and Kimmel 1968).

The Site is relatively flat and underlain by very dense, poorly graded, medium sand and gravel with interbeds of silt to depths of at least 70 ft below ground surface (bgs). These native deposits have been covered by a layer of fill consisting of gravel to silty sands ranging from 1 to 12 ft thick. A low-permeability silt layer has been observed from approximately 44.5 to 47 ft bgs (equivalent to approximate elevations of 200 to 203 ft) in borings advanced in the southwest portion of the Site and is assumed to be a discontinuous low permeability layer within the Vashon outwash. A second low permeability silt layer was also observed from 68 to 70 ft bgs in this area.

1.3.3 Hydrogeology

Groundwater first occurs at depths of approximately 20 ft bgs in an unconfined aquifer in the sandy fluvial deposits. A second regional aquifer lying within older, pre-Vashon deposits underlie the Site and is separated from the shallower glacial outwash aquifer by a semi-confining silt to clay layer observed during investigations at other sites in the vicinity at approximately 100 to 130 ft bgs (URS 2005, Kennedy Jenks 1993).

The Sitewide potentiometric surface and groundwater flow direction have been measured during multiple previous monitoring events. Overall horizontal gradients in the shallow water-bearing zone are relatively flat, with overall variation of less than 1 ft across the Site. Horizontal flow directions are variable, but the overall flow direction appears to be to the west-northwest and influenced by seasonal infiltration of stormwater in the unpaved Sound Transit ROW to the southeast of the property.

Vertical gradients are consistently downward between the uppermost saturated zone and lower saturated zone within the Vashon outwash. Hydrologic studies concerning City of Tacoma Well 12A have found strong downward vertical gradients from the outwash sands to underlying deeper and older glacial deposits, which may influence groundwater flow more strongly than horizontal gradients. The Site lies outside the theoretical capture zone of Well 12A, and only limited pumping effects are observed during periods of groundwater withdrawal at Well 12A (Landau 2024).

Surrounding surface water bodies include Snake Lake approximately 0.7 miles to the southwest and the Thea Foss Waterway arm of Commencement Bay approximately 2.1 miles to the east.

1.3.4 Climate

The climate in the Tacoma area is tempered by winds from the Pacific Ocean, with average daily summer and winter temperatures of approximately 65 degrees and 45 degrees Fahrenheit, respectively. Snow and freezing temperatures are uncommon. Summer rainfall is generally infrequent and light. During the rest of the year, rains are frequent, especially late in fall and winter. The average annual precipitation in the area is approximately 39.8 inches.

2.0 NATURE AND EXTENT OF CONTAMINATION

2.1 Previous Environmental Investigations

Numerous environmental investigations have been undertaken at and adjacent to the Site by various consultants to assess the subsurface impacts from former creosoting operations. Impacts due to former creosote operations were first observed during excavation and grading for the construction of the south warehouse in 1986, when creosote-like material was encountered in surface soil in the vicinity of the former retorts. The following discussion briefly summarizes the investigations. Additional detail is provided in the Draft RI Report (Landau 2024).

- 1) 1992 Phase 2 Environmental Site Assessment. The investigation involved drilling five soil borings (B-01 through B-05) inside the warehouse building and immediately to the east outside of the warehouse and installing three monitoring wells (MW-01, MW-02, and MW-03) near the southeast, northeast, and northwest corners of the warehouse. Soil samples were analyzed for volatile organic compounds (VOCs), gasoline and diesel petroleum hydrocarbons, PAHs, and phenols. Monitoring well groundwater samples were analyzed for VOCs and metals. One groundwater sample collected from MW-02 in the vicinity of the City Materials Testing Lab exceeded the Model Toxics Control Act (MTCA) Method A CUL for trichloroethene (TCE), and all groundwater samples exceeded the MTCA Method A CULs for total chromium and lead. All other soil and groundwater analyte concentrations were less than screening levels.
- 2) 2006 Limited Phase 2 Soil and Groundwater Investigations. The investigation included drilling four soil borings in the former creosoting retort area (SP-01 through SP-04), installing a monitoring well (MW-04) south of the warehouse in the inferred downgradient direction from the former creosoting retorts, and sampling groundwater in MW-01 and MW-04. During soil boring and monitoring well installation, a strong creosote-like odor was noticed in a soil layer with wood fragments encountered between 3 and 6 ft bgs that was interpreted to be the disturbed historical ground surface. Soil and groundwater samples were analyzed for PAHs, including carcinogenic PAHs (cPAHs). All soil samples had cPAH toxic equivalent (expressed as the benzo(a)pyrene toxic equivalent [TEQ]) and naphthalene concentrations exceeding the MTCA Method A CULs. However, total PAHs were less than the MTCA Method A CUL in groundwater samples, and cPAHs were not detected in groundwater.

Five additional soil borings (SP-05 through SP-09) were then drilled outside the warehouse to the south and west, and two additional monitoring wells (MW-05 and MW-06) were installed to the southwest in the inferred downgradient and cross-gradient directions of the former creosoting retort. Six shallow soil samples (SS-01 through SS-06, approximately 1 ft deep) were collected to the east of the warehouse to further investigate the chromium detected in MW-02 groundwater. Soil and groundwater samples were analyzed for PAHs, lead, and chromium. cPAH concentrations exceeded the MTCA Method A industrial-use CULs in soil samples from the three borings immediately west and south of the warehouse (SP-06, SP-07, and SP-08), and cPAH and naphthalene in MW-05 and MW-06 groundwater exceeded MTCA Method A CULs. Chromium and lead were detected at concentrations less than the MTCA Method A CULs in all soil samples.

3) **2008** City of Tacoma Materials Handling Laboratory Remedial Investigation. A remedial investigation (RI) was completed for the former City of Tacoma Materials Handling Laboratory, and it included installation of three permanent monitoring wells (MHLMW-6 through MHLMW-8) screened from approximately 26 to 36 ft bgs and collection of groundwater reconnaissance

samples from temporary well points in three soil borings (HSA-8 through HSA-10, at approximately 25 ft bgs) on the Site. These wells and borings were installed as part of a larger remedial investigation to assess onsite and offsite impacts of prior TCE releases at the Materials Handling Laboratory. Pacific Groundwater Group (PGG) found elevated TCE concentrations in groundwater extending 200 ft onto the northeast corner of the Site, with the extent of the plume delineated by boring HSA-10.

4) **2008 Additional Soil and Groundwater Characterization.** A soil and groundwater investigation was conducted to define the extent of PAHs in soil and groundwater. The investigation included five soil borings (SB-01 through SB-05) to the north, south, east, and west of the former creosoting retort area and three monitoring wells: one (MW-07) located within the footprint of the former creosoting retort area, one (MW-08) located inside the warehouse to the north, and one (MW-09) located adjacent to MW-05 to examine deeper (70 ft bgs) groundwater quality near the former creosoting retorts. Soil samples were analyzed for PAHs, and in one sample (SB 03) collected north of the creosoting operations, cPAH TEQ detections exceeded the MTCA Method A Industrial CUL. Reconnaissance groundwater samples collected from soil borings and monitoring well samples were analyzed for PAHs. Naphthalene concentrations exceeded the MTCA Method A CUL in reconnaissance groundwater samples collected via direct-push temporary well points west of the warehouse (SB-05, MW-07, MW-09). cPAHs exceeded MTCA Method A CULs in all locations except the sample collected along the southeast corner of the warehouse (SB-02). In monitoring well groundwater samples collected from wells to the south and west of the warehouse (MW-05 and MW-06), naphthalene concentrations exceeded MTCA Method A CULs; cPAH concentrations exceeded MTCA Method A CULs in those locations as well as in the warehouse (MW-07) and slightly exceeded MTCA Method A CULs in deeper groundwater (MW-09).

Six additional soil borings (SB-06 through SB-11) were advanced in 2009 within the Sound Transit ROW to the south. Groundwater samples were collected from the temporary wellpoints installed during drilling and analyzed for PAHs including naphthalenes and VOCs. Naphthalene concentrations exceeded MTCA Method A CULs in samples from SB-05 and SB-09, closest to the warehouse. No other samples exceeded MTCA Method A CULs.

- 5) 2014 Focused Subsurface Investigation. A focused subsurface investigation was conducted on the properties on behalf of the owners of the west-adjacent properties in 2014. Nine borings (ECI-B1 through ECI-B9) were drilled to 40 ft bgs for collection of groundwater reconnaissance samples beneath the Shea and Nyssen buildings. At each boring location, two representative soil samples and groundwater samples were collected, one at the water table and one at 40 ft bgs. Samples were analyzed for total petroleum hydrocarbons (TPH), benzene/toluene/ethylbenzene/xylenes (BTEX), and naphthalene. Naphthalene exceeded the MTCA Method A CUL in the deeper groundwater samples at ECI-B1 and ECI-B2 nearest the Shea building property line. Elevated naphthalene concentrations were not detected on the Nyssen property, beneath the Shea warehouse, or in groundwater samples collected at the water table. Concentrations in soil samples did not exceed MTCA Method A CULs, and results in soil were primarily non-detect.
- 6) **2016 Vapor Intrusion Investigation.** Two sub-slab soil vapor samples were collected in 2016, one each in the eastern and western halves of the Shea building. Both samples were analyzed for naphthalene and BTEX. Soil vapor concentrations in both samples were below the MTCA Method B sub-slab soil vapor screening levels.
- 7) **2009–2014 Initial Remedial Investigations.** In June 2010, three monitoring wells were installed off property to further delineate the groundwater naphthalene plume. The additional wells

included MW-10 in the City of Tacoma ROW west of the former creosoting retort area, MW-11 in the ROW to the southwest of the former retort area, and MW-12 to the northwest of the former retort area.

The monitoring well borings were advanced to a depth of 46 ft bgs, with split-spoon samples collected at 5-ft intervals for soil characterization. The upper low-permeability layer, consisting of stiff sandy silt, was encountered at 45 ft bgs at MW-12 but was not reached in the other well borings. The well screens were set immediately above the low-permeability layer from 35 to 45 ft bgs. A slight naphthalene odor was encountered in the lower saturated-zone soil from approximately 35 to 40 ft bgs at MW-10, but a sheen was not observed. Odors and sheens were not encountered at MW-11. A slight naphthalene odor was encountered at MW-12 beginning at 30 ft bgs, and strong odors were noted at 40 ft bgs; a sheen was observed on groundwater on the split-spoon sampler but was not replicated during a sheen test performed on soil from the MW-12 interval with strong odor. Soil samples were collected from the lower saturated interval where creosote odors were noted but ultimately were not analyzed because of the absence of sheen, non-aqueous phase liquids (NAPLs), or other indications of the presence of source contamination in soil.

One soil boring (SP-2B) was also installed within the warehouse in the vicinity of the former creosoting retort area in June 2010. Intermittent oily odor and trace sheen were observed at SP-2B beginning approximately at 8 ft bgs, and a strong naphthalene odor and sheen were observed below approximately 30 ft bgs. Soil samples were collected at 15, 30, 35, and 45 ft bgs and analyzed for PAHs and naphthalene. The maximum concentration of naphthalene measured in soil was 470 milligrams per kilogram (mg/kg) at 30 ft bgs. The maximum concentration of cPAHs was 9.2 mg/kg, also at 30 ft bgs.

The existing monitoring well network was sampled in June 2010 and December 2011, with groundwater samples from all wells analyzed for PAHs and from a subset of wells for VOCs, including naphthalene. The maximum naphthalene concentration detected during the two sampling events was 14,000 micrograms per liter (μ g/L) at MW-10 in December 2011. The maximum concentration of cPAHs was 8.2 μ g/L detected at MW-06. Elevated concentrations of benzene, 1,2,4-methylbenzene, 1,3,5-methylbenzene, ethylbenzene, and xylenes were also detected in MW-10 and MW-12.

Quarterly groundwater monitoring was conducted at a subset of Site monitoring wells from 2012 to 2013 to confirm that naphthalene concentrations were stable within a reasonable range of seasonal fluctuations. Quarterly monitoring samples were collected in November 2012 and March, June, and October 2013. The quarterly monitoring network included MW-01, MW-05, MW-06, MW-07, MW-10, MW-11, and MW-12. The deeper aquifer well MW-09 was also sampled during the first and last quarterly events. Results from the 2012–2013 quarterly sampling were generally consistent with the 2010–2011 sampling events, with significantly elevated naphthalene concentrations in MW-06, MW-10, and MW-12, and a maximum naphthalene concentration of 14,000 $\mu g/L$ detected in MW-10. Concentrations of ethylbenzene and xylenes were also detected at elevated concentrations in MW-10 and MW-12.

In 2011 and 2012, temporary well points (SB-21 through SB-29, SB-31 through SB-39) were installed to delineate the extent of the naphthalene plume in groundwater, with groundwater reconnaissance samples analyzed for VOCs, including naphthalene. Naphthalene detections in these groundwater reconnaissance samples ranged from non-detect in samples collected to the north, south, and east of the property to a maximum of 17,000 μ g/L in the sample collected at 40 ft bgs from the MW-9 boring directly west of the former creosoting retort area. Borings to

the west and northwest of the former creosoting retort area (SB-29, SB-21, SB-23, and SB-31 through SB-36) also yielded elevated naphthene concentrations, with a maximum concentration of $5,400 \mu g/L$ in SB-29.

In December 2011, indoor air samples for naphthalene analysis were collected from three locations inside the warehouse building and one upwind outdoor area immediately after the building had been shut for the weekend. The naphthalene concentrations measured in indoor air ranged from 0.82 to 1.1 micrograms per cubic meter ($\mu g/m^3$), less than the applicable MTCA Method B CUL of 1.4 $\mu g/m^3$ at the time of sampling.

8) 2018–2021 Supplemental Remedial Investigations. In 2018, one monitoring well (MW-14) and four soil borings (SB-40 through SB-43) were installed to delineate the extent of groundwater contamination in and around the former retort area. The boring for MW-14 was advanced to 80 ft bgs, and based on field-screening, the well screen was set within the most impacted shallow groundwater interval from 35 to 45 ft bgs. Soil samples were collected from the MW-14 boring at the observed creosote-impacted interval and at three underlying intervals and analyzed for VOCs and cPAHs. Naphthalene and cPAHs were detected at concentrations of 167 mg/kg and 8.45 mg/kg, respectively, in a soil sample collected from 20.5 to 21 ft. Soil borings SB-40 through SB-43 were advanced to depths of 80 to 85 ft bgs, with the upper low-permeability silt layer observed in all borings at depths of 47 to 49 ft bgs (measured from the warehouse floor grade) and the lower low-permeability silt layer observed in all borings except SB-41 at depths ranging from 69 to 72 ft bgs. The soil most heavily impacted by creosote was observed at SB-43 in the western portion of the former creosote retort. At this location, strong odor and rainbow sheen were observed between approximately 25 and 40 ft bgs. Creosote impacts were also noted in SB-40 (approximately 27 to 40 ft bgs) and in SB-41 and SB-42 (approximately 25 to 28 ft bgs). Representative soil samples were collected from the impacted intervals and underlying intervals without field indications of contamination at each boring location. Samples were analyzed for VOCs and cPAHs, and a representative highly contaminated sample from SB-43 was additionally analyzed for pentachlorophenol.

In January 2019, reconnaissance groundwater samples were collected at the targeted locations for new wells MW-15, MW-16, and MW-17. These samples were collected for screening purposes prior to permanent well installation, to refine the intended monitoring locations at the downgradient edges of the groundwater contaminant plume. Based on field observations of naphthalene odor and screening sample results with elevated naphthalene, the permanent well locations at MW-16 and MW-17 were both shifted northward. In August 2020, MW-15, MW-16, and MW-17 were installed to fully delineate the groundwater naphthalene plume. The wells were installed inside the west-adjacent Shea warehouse (MW-15), outside the northwest-adjacent Nyssen building (MW-16), and on the property to the north of MW-12 (MW-17). The monitoring well borings were advanced to depths of 45 to 50 ft bgs; the upper low-permeability layer was encountered at depths of 41.5 to 43.5 ft bgs, and the well screens were set immediately above it. Slight naphthalene odor was encountered only in saturated-zone soil from approximately 39 to 40 ft bgs at MW-16, and sheens were not encountered in any well boring locations. Soil samples for laboratory analysis were not collected because soil was not observed to be impacted.

An additional four quarters of monitoring data were collected from the Site monitoring well network to fully delineate the groundwater naphthalene plume in September and December 2020, and in March and June 2021. The well network included MW-01, MW-05, MW-06, MW-09, MW-10, MW-11, MW-12, MW-14, MW-15, MW-16, and MW-17. MW-04 and

MW-07 were not sampled because their screened intervals were not sufficiently deep to sample the most contaminated zone of groundwater immediately above the uppermost low-permeability layer, and MW-01 was not sampled because of inaccessibility and location relative to the former creosoting retort area. During the first quarter of monitoring, samples were analyzed for VOCs, and selected samples from new wells (MW-14 through MW-17) and wells near the former creosoting retort area (MW-05, MW-06, and MW-09) were additionally analyzed for cPAHs and other selected semivolatile organic compounds (SVOCs). Based on the results of the first round of quarterly monitoring, during which only samples collected from MW-06 had detectable cPAHs, cPAHs and other SVOCs were monitored only at MW-06. MW-05 has a shallower screened interval than other shallow wells in the monitoring network, and no contaminants were detected in samples collected from it during the first round of sampling despite MW-05 being located adjacent to the former creosoting retort. It was determined that MW-05 was also screened too shallow to sample the most contaminated zone of groundwater, and it was subsequently removed from the monitoring network. The monitoring well network for the remaining three quarterly monitoring events in 2020–2021 consisted of shallow wells MW-06, MW-10, MW-11, MW-12, MW-14, MW-15, MW-16, and MW-17, and deep well MW-09.

To assess the potential for vapor intrusion, two sub-slab soil vapor samples (VP-01 and VP-02) were collected within the south warehouse in 2020. The vapor points were installed adjacent to the locations of existing soil borings SP-2 and MW-07, where previous field-screening identified the strongest indications of creosote contamination in shallow soil. Sub-slab vapor sample points VP-01 and VP-02 were installed by drilling through the warehouse floor with a roto-hammer and installing a vapor pin implant sealed with a silicone gasket. The sample points were allowed to equilibrate for a minimum of 1 week after installation prior to sample collection.

2.2 Summary of Results

Appendix A provides the logs for soil borings and monitoring wells installed at the Site. Figure 1 summarizes the RI soil and groundwater CUL exceedances, and RI Figures 5.1 through 5.3 (see Appendix B) provide more detailed depictions of the Site data. The RI tables (see Appendix C) provide tabulated summaries of the proposed Site CULs and data.

Soil contamination is generally limited to the area beneath and immediately adjacent to the former creosoting retort area, except for limited exceedances of the proposed CUL for cPAHs to the north (which are likely associated with historic regrading). The vertical extent of contamination in soil and groundwater is bounded by the lowermost extent of a silt confining layer from 45 to 50 ft bgs, below which there are no exceedances of CULs for naphthalene or cPAHs in soil.

The naphthalene and benzene/xylene groundwater plume radiates in all directions from the presumed source area in the vicinity of the former creosoting retort. While the delineated plume shows no strong flow direction, the overall greatest migration aligns with the horizontal hydraulic flow direction to the north-northwest. Due to the low hydraulic gradient (typically less than 1 ft Site-wide) and biodegradation, the plume has largely been static over time with slight seasonal variability.

2.3 Conceptual Site Model

This section provides a summary of the conceptual site model (CSM) for contamination related to the former creosote plant and is discussed in more detail in the draft RI report. The CSM demonstrates the current understanding of the nature and extent of contamination relative to the presumed historic source area in the vicinity of the former creosoting retort and migration pathways to receptors.

2.3.1 Potential Sources of Contamination

The loading and removal of creosote or pipe from the early 1900s to the mid-1930s would likely have caused small to moderate releases of creosote. During historical Site operations, creosote releases that likely originated as surface spills were able to saturate soil and migrate downward through the unpaved, permeable ground surface and factory floors. Releases were most likely near or under the approximate footprint of the historical creosoting retorts (shown on Figure 2) and are the cause of present-day contamination on the Site. A significant portion of the creosote-contaminated soil in the vicinity of the former creosote retort was removed during construction of the warehouse that presently occupies the Site. The creosote soil footprint is currently below the southern portion of the warehouse. There are no continuing sources of hazardous substances stored or used at the Site.

Storage tanks containing gasoline were removed from the Site in 1989. Prior to the removal of gasoline storage tanks in 1989, no gasoline releases were documented onsite, and contamination related to potential releases have not been identified during subsequent environmental investigations.

2.3.2 Contaminant Transport Pathways

Based on the current understanding of the Site, current land use, and numerous environmental investigations, three primary transport mechanisms have been identified:

1) Historical gravity-driven vertical migration of NAPL creosote within the vadose and saturated zones. The primary transport mechanism in vadose-zone soil was the gravity-driven downward migration of relatively dense NAPL creosote from releases during historical Site operations from the early 1900s to the mid-1930s. The inherent viscosity of creosote tends to limit its transport in soil, and cPAHs and naphthalene tend to adsorb strongly to organic material in soil, further limiting their transport. However, Site investigations have shown that creosote likely slowly migrated downward through the vadose zone and saturated zone over the last 90 years due to its relative density. Downward migration of the creosote continued to the silt confining layer at 45 ft bgs, with a significant portion of the soil contamination at the Site currently located between depths of approximately 20 and 45 ft bgs. The source area is defined by the presence of creosote product, including visible creosote stringers and sheen-producing soil, that occurs primarily in the 30 to 45 ft bgs depth interval in the vicinity of the former creosote retort. Within this source area, the greatest concentrations of naphthalene, benzene, and xylenes in groundwater are located within the approximate 35 to 45 ft bgs interval directly above the low permeability silt layer. Markedly elevated contaminant concentrations are present in samples collected from this interval versus the overlying portions of the saturated zone. The stratification of contamination in the upper saturated zone is attributed to biological degradation by aerobic bacteria in the more oxygen-rich upper interval, compared to a lesser

- rate of biodegradation in lowermost interval. Strong downward hydraulic gradients may additionally influence the distribution of contaminants in the upper saturated zone.
- 2) Leaching of soluble fractions of the creosote mixture in the saturated zone and limited migration downward and laterally in the dissolved phase. The Site has been paved for its recent operational history, thereby preventing stormwater infiltration and preventing leaching of contaminants from vadose-zone soil into the shallow aguifer. Therefore, the primary mechanism driving ongoing contaminant migration in groundwater is the leaching of soluble naphthalene and benzene from creosote-contaminated source soil in the saturated zone. Beyond the area of residual soil contamination in the source area, the extent of naphthalene in groundwater appears to be dictated by the flat hydraulic gradient to the north-northwest onsite and by localized secondary flow directions. A "halo" of elevated naphthalene concentrations in groundwater beyond the boundary of the source area is present in the upgradient and crossgradient directions due to these secondary flow directions rather than to the presence of contaminated soil in those areas. Groundwater data from the borings surrounding the monitoring well network demonstrate that the elevated naphthalene concentrations in the vicinity of the former creosoting retort area decreases substantially with increased distance from the former retort. This decrease is attributable to the degradation of naphthalene (particularly at the edges of the plume).
- 3) Volatilization of naphthalene, benzene, and xylene to soil gas in the vadose zone. Chemicals associated with historic creosoting operations are volatile and semivolatile in nature. These chemicals present a potential risk to indoor air quality if present in high concentrations and if structures are located or built over contaminated areas.

2.3.3 Potential Receptors and Exposure Pathways

The three primary exposure pathways of concern at the Site are direct contact with groundwater via consumption of drinking water, worker direct contact with soil, and soil leaching to groundwater:

- Consumption of Drinking Water. Groundwater at the Site is not currently used for drinking water but is potentially potable. Therefore, the exposure pathway of direct contact via consumption of drinking water is considered complete at the Site.
- 2) Direct Contact. Impacted soil at the Site is not currently accessible via direct contact due to the presence of overlying buildings or pavement but could be contacted by workers or other receptors if future development were to occur. The soil direct contact pathway is considered incomplete for terrestrial ecological receptors at the Site. Leaching from impacted soil to groundwater via infiltration of surface water likely occurred in the past before construction of the current impervious surfaces of buildings and pavement in the former creosoting retort area. Leaching from impacted soil due to groundwater flow through impacted soil in the saturated zone to groundwater is presumed to be ongoing.
- 3) Inhalation of Indoor Air. The results of sub-slab soil vapor samples collected during the RI indicate that the vapor intrusion pathway is not complete under current conditions. The pathway may need to be reevaluated if conditions change (e.g., redevelopment).

The RI determined that due to the type and depth of Site contamination, the paved surface of the Site, and the location of the nearest residential area, there is a low risk of environmental exposures to residential communities, vulnerable populations, and overburdened communities from this Site.

3.0 PILOT STUDY APPROACH

This section discusses the approach for conducting a pilot study at the Site, including the objectives of the study, the technologies considered, and the general description of the pilot study to be conducted using the selected technology. Section 4.0 provides the pilot study implementation details.

3.1.1 Pilot Study Objectives

The objectives of the pilot study are to (1) apply a technology to reduce contaminant mass, contain contaminant source material, and lower contaminant flux in the creosote source area, (2) evaluate the results of the application, (3) provide a recommendation for potential expanded use of the technology in the source area, and (4) provide information to complete an appropriate feasibility study (FS) for the Site.

3.1.2 Technology Evaluation and Selection

While multiple technologies are available for remediating creosote source areas, Landau identified three priority technologies for use in a source-area pilot study, each potentially applicable for reducing creosote source-area mass and contaminant flux and each appearing to have the highest likelihood for good value-to-cost ratios. Landau additionally evaluated several other technologies (e.g., thermal treatment technologies, *in situ* solidification/stabilization, and excavation and removal) but did not prioritize those for the pilot study given the likely low value-to-cost ratio or limited applicability due to Site constraints. Following is a brief discussion of each of the prioritized technologies:

- Enhanced In Situ Bioremediation (EISB). This broad class of treatment typically involves injecting a combination of an organic substrate, nutrients, other amendments, and/or specialized microbes to facilitate the biodegradation of contaminants. EISB can occur aerobically or anaerobically depending on the contaminant to be treated. This type of injection creates conditions that stimulate bacterial growth and, since the injected materials are soluble in water, may flow some distance downgradient from the point of injection. While this technology can treat contaminants sorbed to soil and dissolved in pore water with relatively minimal site disturbance, Landau has not selected this technology for use in the creosote source area pilot study due to the decreased availability for biodegradation of high-concentration creosote source material (e.g., higher ring PAHs) and the correspondingly limited efficiency of the treatment in high-mass areas.
- In Situ Chemical Oxidation (ISCO). This type of treatment involves the introduction of a chemical oxidant into the subsurface for the purpose of transforming groundwater or soil contaminants into less harmful chemical species. Examples of oxidants include sodium permanganate, sodium persulfate, ozone, and Fenton's Reagent (a hydrogen peroxide/ferrous ion solution). In general, oxidants will persist in the subsurface up to several months for permanganate and persulfate applications and may require reapplication to enhance effectiveness. This technology destroys contaminant mass in situ relatively rapidly and with minimal site disturbance.
- In Situ Geochemical Stabilization (ISGS). Source area treatment by ISGS involves an in situ injection of an ISCO chemical (sodium permanganate) along with binders and reagents designed to both treat high-concentration source material (NAPL or dissolved phase) through oxidation as

well as reduce the potential for contaminant leaching of the residual NAPL and/or high-concentration soil. The ISGS process deposits precipitates around the NAPL, hardening and chemically weathering the NAPL and increasing the NAPL viscosity, which reduces both the flux of contaminants from the NAPL and the hydraulic conductivity of the aquifer in the target treatment zone. This technology destroys contaminant mass *in situ* in a relatively rapid timeframe, stabilizes NAPL in the source area, requires relatively simple materials and equipment, generates no significant waste, and minimizes site disturbance.

Based on the pilot study objectives and the location of the source area within the warehouse, Landau has selected the ISGS technology for use in the pilot study. It both contains and treats the source mass, reduces flux from the source mass to groundwater, and reduces the hydraulic conductivity of the aquifer in the target treatment zone with relatively minimal site disturbance, exhibiting some of the positive aspects of both ISCO and *in situ* solidification/stabilization technologies. The ISGS technology has been successfully employed in other creosote source areas around the country, and Landau expects the technology to be productive at the Site.

3.1.3 Pilot Study Description

The pilot study will consist of a single injection of ISGS reagents within what appears to be the primary dense nonaqueous phase liquid (DNAPL) horizon in the western portion of the former creosoting retort area, monitoring the effects of the ISGS injections, evaluating the results, and preparing a report documenting the pilot study.

The ISGS reagents will be injected in new polyvinyl chloride (PVC) injection wells installed in the western portion of the former creosote source area in the vicinity of boring SB-43, which had strong evidence of creosote contamination during drilling and the highest soil naphthalene in the source area. Section 4.3 discusses the layout and preliminary design of the injection wells. Based on observations made during injection well installation and baseline soil and groundwater sampling conducted in the injection wells, the layout and design of the injection and performance monitoring networks may be modified.

Performance monitoring will be conducted in two existing monitoring wells and in five new monitoring wells installed around the perimeter of the target treatment zone. Monitoring will take place before, during, and after the injection event (see Section 4.5).

The ISGS product to be used in the pilot study will be Evonik's ISGS technology, which consists of sodium permanganate and a proprietary blend of cold-water soluble inorganic metals and salts. The technology is licensed to Innovative Environmental Technologies, Inc. (IET), which provides the product and conducts the injections. Landau will set up the pilot study, coordinate schedule and field activities with stakeholders, coordinate the installation of the injection wells and monitoring wells, develop the wells, coordinate the injection event, provide oversight during fieldwork, conduct the performance monitoring, and incorporate the pilot study results in the FS.

4.0 PILOT STUDY IMPLEMENTATION

This section describes the field activities related to pilot study implementation, including treatability testing, well installation, amendment injection, groundwater monitoring, equipment decontamination, and management of residual wastes. Field activities will be conducted in accordance with a Site-specific Health and Safety Plan (HASP), which will be prepared prior to initiation of field activities.

4.1 Cultural Resource Review

Prior to ground-disturbing activities during the implementation of the pilot study work plan and consistent with Washington Administrative Code (WAC) 173-340-815(3)(a), Landau will initiate a cultural resource review by Ecology. An Inadvertent Discovery Plan will be prepared to avoid, minimize, or mitigate the effects of the pilot study field activities on archaeological and cultural resources.

4.2 Underground Injection Control Registration

Prior to the pilot study injection, the Site will be registered with Ecology's Underground Injection Control (UIC) program per Chapter 173-218 of the WAC. Per WAC 173-218-040, the injection wells at the Site are considered Class V injection wells because they will receive "fluids intended to clean up, treat or prevent subsurface contamination." Class V UIC wells must be registered and either rule-authorized (WAC 173-218-070) or receive a state waste discharge permit issued by Ecology.

4.3 Injection and Monitoring Well Installation

Landau will subcontract a state-licensed well driller to install seven injection wells and five monitoring wells. Figure 3 shows the approximate locations of the wells, which may be moved based on the presence of utilities or due to warehouse operational needs. The driller will install the wells with either a hollow-stem auger or sonic drill rig. Based on the observations during drilling of soil borings and monitoring wells in the former creosote retort area and the distribution of naphthalene and PAHs in soil and groundwater, the injection wells will be installed with screens between 25 ft and 35 ft bgs (which is the warehouse floor; up to 3 ft above the surrounding grade; see Table 1). Based on the relatively coarse and likely permeable nature of the aguifer, it is anticipated that the radius of influence (ROI) in the injection wells will be at least 10 ft. Allowing for 1 ft of overlap in ROIs, the injection wells will be spaced approximately 19 ft apart (see Figure 3). The injection wells will be installed, developed, and sampled before installation of the performance monitoring wells. Based on observations made during drilling and the results of the baseline groundwater sampling conducted in the injection wells and existing monitoring wells, the layout and design of the injection network and performance monitoring network may be modified. For example, Figure 3 shows the location of two potential contingency injection well locations that may be installed and used based on field observations of an abundance of DNAPL in the injection well borings on the west side of the target treatment zone or elevated baseline groundwater sampling results in existing monitoring wells MW-05 and MW-09.

Five performance monitoring wells will be installed around the target treatment zone. All monitoring wells will be screened between 25 ft and 35 ft bgs (relative to the warehouse floor). Two wells will be

installed on the north side of the target treatment zone, two wells will be installed on the south side of the target treatment zone, and one well will be installed on the east side of the target treatment zone, with all wells located approximately 10 to 15 ft outside the target treatment zone. Because existing monitoring wells MW-05 (screened 25 to 35 ft bgs) and MW-09 (screened 60 to 70 ft bgs) are located to the west of the former creosote retort area, no additional performance monitoring wells will be installed on the west side of the target treatment zone.

The wells will be constructed of Schedule 40 PVC well materials, with silica sand backfill, a bentonite annular seal, and a flush-mount surface monument set in concrete. Landau will develop each monitoring well using surging, bailing, and pumping techniques to clean out silty water introduced during the well installation process, establish a hydraulic connection between the aquifer and the well filter pack, and reduce the groundwater turbidity in the well.

A state-licensed surveyor will determine the horizontal and vertical coordinates of the wells. The elevations of the wells will be relative to NAVD88, and the horizontal locations of the wells will be relative to NAD83/91,² Washington State Plane Coordinate System North Zone.

4.4 In Situ Geochemical Stabilization Reagent Injection

This section provides the general procedures to be followed during the ISGS reagent injections. The procedures may be refined or modified based on results of the initial injection well installation and treatability study.

4.4.1 Pilot Study Volume and Dosing

The preliminary design volume for the pilot study injections at the Site will be based on the following:

- ROI = 10 ft
- ROI overlap = 1 ft
- Treatment area = 32 ft by 72 ft, or approximately 2,300 SF
- Treatment length per well = 10 ft
- Effective porosity = 20 percent
- Target percentage of pore space to fill with ISGS reagent solution = 11 percent.

Based on these parameters and the conservative upper end of the typical range in reagent injection percentages (10 percent reagent), approximately 49,100 pounds of ISGS reagent will be injected during the pilot study.

4.4.2 In Situ Geochemical Stabilization Reagent Material Compatibility

Sodium permanganate is compatible with plastics (including polyethylene, polypropylene, and Teflon), many metals (including carbon steel and 316 stainless steel in neutral and alkaline solutions), and synthetic materials. Sodium permanganate is often incompatible with natural rubbers and fibers, so

² North American Datum of 1983/1991.

contact with those materials should be avoided. Aluminum, zinc, copper, lead, and alloys containing these metals may be (slightly) affected by sodium permanganate solutions.

4.4.3 Mixing and Injection Procedures

Due to corrosivity, staff will wear compatible personal protective equipment during mixing and injection. Safety glasses, gloves, and PVC splash bibs or aprons will be worn. Health and safety measures will be detailed in the forthcoming HASP.

Sodium permanganate will be shipped to the Site in 3,000-pound caged plastic totes. Generally, to create the injection solution, the permanganate will be mixed with tap water and the other amendments in batches. Portable tanks and motorized paddles located in an injection trailer will be used to mix the solution. The solution will be pumped using a compressed, air-powered, double-diaphragm injection pump (made from compatible materials) into each injection well through a flexible chemically compatible hose (with a minimum pressure rating of 250 pounds per square inch [psi]) affixed to a cam lock at the wellhead. Pumping will continue until the design volume has been pumped into the well. Injection pressure and volume will be monitored throughout the process. At the end of the injection at each well, clean water will be pumped into the well to clean it out for potential future use.

More specifically, IET will use the following procedures and equipment.

4.4.3.1 Procedures

- 1) **Subsurface Pathway Development.** Initially, compressed air will be delivered to the subsurface via IET's proprietary injection trailer system. This process allows for confirmation of open delivery routes while enhancing horizontal injection pathways. By confirming open and viable subsurface delivery pathways, it is anticipated that upon introduction of the ISGS reagent, the injectate will flow freely, thus minimizing health and safety risks associated with full oxidant injection lines and injection tooling when no subsurface delivery route has been established. Confirmation of open and free pathways is accomplished via observed pressure drops and free moving compressed gases to the subsurface.
- 2) **ISGS Emplacement.** A 10 percent solution of ISGS will be introduced at pressures between 15 and 50 psi and flow rates between 15 and 25 gallons per minute (gpm). A small amount of water will then be introduced to rinse the injection equipment.
- 3) Post Liquid injection Compressed Air Injection. The injection lines will be cleared of liquids, and all injectants will be forced into the created formation and upward into the vadose zone. This step is performed so that all material will be injected outward into the formation, minimizing any surface excursions of injectants following the release of the injection pressure. Once the injection cycle is complete, the injection point will be temporarily capped to allow the pressurized subsurface to accept the injectants.

Spill Prevention. Spill prevention measures will be used during injection activities to prevent the accidental release of injection fluid. Mixing tanks, injection pumps, and major hose connections will be placed in containment berms. Pumps and hoses will be attended and monitored during injection. All hose connections will be secured with zip ties. A wet-dry vacuum will be kept onsite during injection activities to immediately collect spilled fluid.

4.4.3.2 Equipment

The injections will occur via IET's mobile oxidation injection trailer as described:

- Injection Lines. High-pressure stainless steel, braided rubber, 1-inch diameter hoses.
- Injection Trailer (see photo below). IET self-contained injection trailer, consisting of two 200-gallon conical tanks capable of maintaining 30 percent solids as a suspension via lightning mixers; onboard generator, all stainless-steel piping system, 2-inch pneumatic diaphragm pump with an operating pressure of 110 psi.; onboard 37 CFM/175 pounds per square inch gauge (psig) compressor with 240 gallons of air storage; and self-contained eye wash and safety shower.

IET INJECTION SYSTEM UNITED STATES PATENT 7,044,152



Injection Trailers Include: Multiple Liquid Feed Systems, Stainless Steel Piping, Isolated Compressed Gas Containment, Safety Shower, Eyewash Station, Onboard Generator, Chemical Resistant Construction, Mobile Office Space



4.5 Performance Monitoring

Performance monitoring will be conducted to evaluate the short-term effects of amendment injections on groundwater levels and groundwater flow directions near the pilot study treatment zone and longer-term effects of amendment injections on groundwater quality near the pilot study treatment zone.

Table 2 summarizes the performance monitoring well network and monitoring frequency.

4.5.1 Water Level Monitoring

Water-level elevations will be monitored in the five new performance monitoring wells (MW-18 to MW-22) before, during, and after pilot study injection activities. Baseline water levels will be measured prior to amendment injection, and water levels will be measured periodically during the injection period, immediately after injection activities have been completed, and during quarterly groundwater sampling events. Water-level elevations will also be monitored in accessible existing Site monitoring wells and in the injection wells during the baseline and quarterly groundwater sampling events.

4.5.2 Groundwater Sampling

Performance groundwater sampling will consist of one baseline sampling event prior to amendment injection and quarterly groundwater sampling beginning approximately 3 months after amendment injection to evaluate treatment effectiveness. During each monitoring event, samples will be collected using low-flow sampling procedures from two existing monitoring wells (MW-05 and MW-09) and five new performance monitoring wells (MW-18 to MW-22). During the baseline sampling event, the injection wells (IW-01 through IW-07) will also be sampled, and accessible existing monitoring wells (MW-04, MW-06, MW-07, MW-10, MW-11, MW-12, MW-14, MW-15, MW-16, and MW-17) will also be sampled during the baseline and final quarterly events. Field parameters (dissolved oxygen [DO], oxidation-reduction potential, conductivity, temperature, and pH) will be measured during well purging, groundwater samples will be submitted to an Ecology-accredited laboratory for analysis of BTEX and naphthalene by US Environmental Protection Agency (EPA) Method 8260D and PAHs by EPA Method 8270E with selected ion monitoring (SIM). One trip blank and one duplicate sample will be analyzed during each sampling event.

4.6 Equipment Decontamination

All non-dedicated sampling and injection equipment will be decontaminated according to the procedures described in the forthcoming HASP. This includes water-level monitoring instruments, sample tubing, and any other equipment that comes into contact with Site soil or groundwater. Equipment will be dedicated, when possible, to reduce the potential for cross-contamination.

4.7 Residual Waste Management

Investigation-derived waste (IDW) will consist of soil generated during drilling activities, water used for decontamination, rinsate from tanks and drums used for mixing injection solution, purge water from sampling of monitoring wells, concrete cores removed during well installation, and personal protective equipment (PPE). Concrete cores and PPE will be handled as municipal solid waste. Residual soil and

Pilot Study Work Plan West Coast Door Site–Tacoma, Washington

water will be stored either in drums or waste bins prior to characterization, profiling, and transportation for disposal at an appropriate offsite facility.

5.0 DATA EVALUATION AND REPORTING

Groundwater monitoring data collected after the pilot study injection will be used to evaluate the effectiveness of the ISGS technology, the need for additional source-area treatment, and full-scale injection design, if necessary. Data will also be used to determine the need and appropriate timing for active dissolved plume treatment and or monitored natural attenuation. After receiving the results of the four quarterly performance monitoring events, Landau will prepare a pilot study completion report documenting the treatability study, well installation and development, ISGS reagent injection, and performance monitoring. The report will include a summary of the installation process; a property map updated with the locations of the injection and performance monitoring wells; tables summarizing the survey results, well completion details, groundwater levels, and analytical data; groundwater elevation contour maps, as needed; well logs; and laboratory analytical reports.

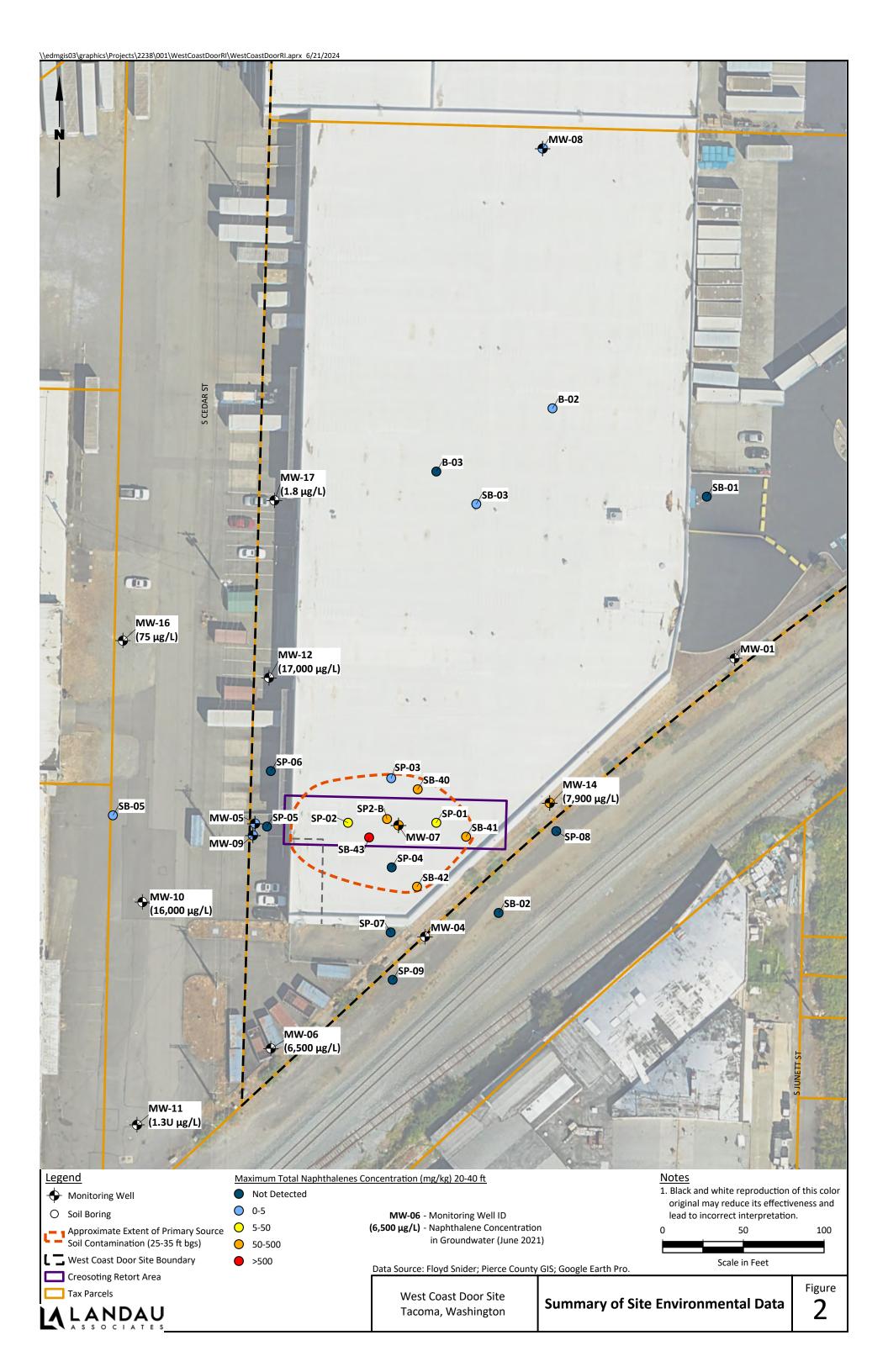
6.0 USE OF THIS WORK PLAN

This work plan has been prepared for the exclusive use of 3102 TIC and their associated legal and insurance carrier teams for the pilot study at the former West Coast Door Facility in Tacoma, Washington. No other party is entitled to rely on the information, conclusions, and recommendations included in this document without the express written consent of Landau. Further, the reuse of information, conclusions, and recommendations provided herein for extensions of the project or for any other project, without review and authorization by Landau, shall be at the user's sole risk. Landau warrants that within the limitations of scope, schedule, and budget, our services have been provided in a manner consistent with that level of care and skill ordinarily exercised by members of the profession currently practicing in the same locality under similar conditions as this project. Landau makes no other warranty, either express or implied.

7.0 REFERENCES

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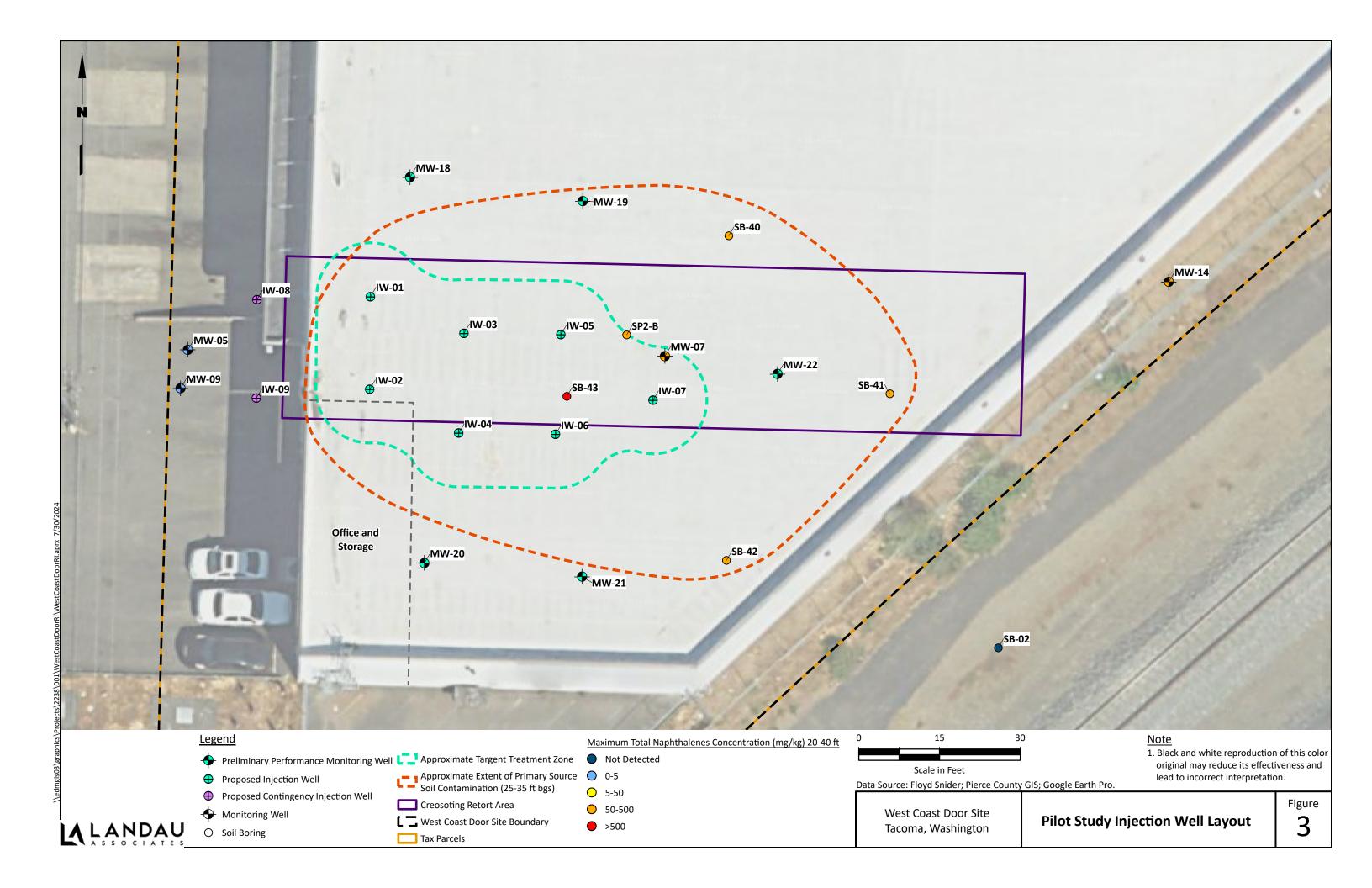


Table 1
Site Well Completions
West Coast Door Site, Tacoma, Washington

Well ID	Installation Date	Total Depth (ft bgs)	Screened Interval (ft bgs)	Top of Well Casing Elevation (ft NAVD 88)	Surface Completion	Notes
Existing Wells		, 0,	, ,,	,	·	
MW-01	6/18/1992	44	23.5–43.5	247.02 ⁽¹⁾	Flush-mounted	Missing; under pavement
MW-02	6/18/1992	44 ⁽²⁾	23.5–43.5 ⁽²⁾	247.36	Flush-mounted	
MW-03	6/18/1992	45	23.5-43.5	247.55	Flush-mounted	
MW-04	7/28/2006	35	25–35	246.27	Flush-mounted	
MW-05	9/16/2006	35	25–35	246.09	Flush-mounted	
MW-06	9/18/2006	35	25–35	245.97	Flush-mounted	
MW-07	1/26/2007	40	25–40	248.18	Flush-mounted	
MW-08	1/31/2007	40	25–40	248.24 ⁽¹⁾	Flush-mounted	Missing; under warehouse partition wall
MW-09	9/7/2007	70	60–70	245.99	Flush-mounted	
MW-10	6/10/2010	46	35–45	244.22	Flush-mounted	
MW-11	6/10/2010	46	35–45	243.55	Flush-mounted	
MW-12	6/11/2010	46	35–45	243.97	Flush-mounted	
MW-14	10/5/2018	45	35–45	250.13	Flush-mounted	
MW-15	8/11/2020	43.5	33.5–43.5	248.28	Flush-mounted	
MW-16	8/10/2020	41	31–41	246.97	Flush-mounted	
MW-17	8/10/2020	41.5	31.5–41.5	247.91	Flush-mounted	
Proposed Inje	ection Wells					
IW-01	-	35	25–35	_	Flush-mounted	Projected depths relative to warehouse floor
IW-02	-	35	25–35	_	Flush-mounted	Projected depths relative to warehouse floor
IW-03	-	35	25–35	_	Flush-mounted	Projected depths relative to warehouse floor
IW-04	-	35	25–35	_	Flush-mounted	Projected depths relative to warehouse floor
IW-05	-	35	25–35	_	Flush-mounted	Projected depths relative to warehouse floor
IW-06	-	35	25–35	_	Flush-mounted	Projected depths relative to warehouse floor
IW-07	-	35	25–35	-	Flush-mounted	Projected depths relative to warehouse floor
Proposed Cor	ntingency Injectio	n Wells				
IW-08	_	35	25–35	-	Flush-mounted	Projected depths relative to ground surface
IW-09	-	35	25–35	_	Flush-mounted	Projected depths relative to ground surface

Table 1
Site Well Completions
West Coast Door Site, Tacoma, Washington

Well ID		Total Depth (ft bgs)	Screened Interval Top of Well Casing (ft bgs) Elevation (ft NAVD 88)		Surface Completion	Notes					
Proposed Per	Proposed Performance Monitoring Wells										
MW-18	_	35	25–35	ı	Flush-mounted	Projected depths relative to warehouse floor					
MW-19	-	35	25–35	-	Flush-mounted	Projected depths relative to warehouse floor					
MW-20	-	35	25–35	-	Flush-mounted	Projected depths relative to warehouse floor					
MW-21	_	35	25–35	- 1	Flush-mounted	Projected depths relative to warehouse floor					
MW-22	_	35	25–35	_	Flush-mounted	Projected depths relative to warehouse floor					

Notes:

- (1) Well not accessible during 2020 re-survey; casing elevation is estimated from prior surveyed elevation in NGVD 29 using a local correction factor of +3.38 ft.
- (2) Well installation log not available, estimated from concurrent MW-01 and MW-03 construction details.

For existing wells, the table has been modified from Floyd | Snider 2022 Draft RI/FS Report.

Abbreviations:

bgs below ground surface

ft fee

NAVD 88 North American Vertical Datum of 1988 NGVD 29 National Geodetic Vertical Datum of 1929

Table 2
Pilot Study Performance Monitoring
West Coast Door Site, Tacoma, Washington

		Post Injection Monitoring					
Well ID	Baseline Sampling	First Quarter	Second Quarter	Third Quarter	Fourth Quarter	Notes	
Existing V		Qualite.	Quai tei	Quai tei	Quarter		
MW-04	X				Х	Sampled if accessible	
MW-05	X	Х	X	X	X	Water level monitoring during injection event	
MW-06	X				X	water level monitoring during injection event	
MW-07	X				X		
MW-09	X	X	X	X	X	Water level monitoring during injection event	
MW-10	Х				Х		
MW-11	Х				Х		
MW-12	Х				Х		
MW-14	Х				Х		
MW-15	Х				Х	Located in the Shea Property warehouse to the west	
MW-16	Х				Х		
MW-17	Х				Х		
Proposed	Injection W	/ells			•		
IW-01	Х						
IW-02	Х						
IW-03	Х						
IW-04	Х						
IW-05	Х						
IW-06	Х						
IW-07	Х						
Proposed	Contingend	y Injection	Wells				
IW-08						Sampled during baseline event if installed	
IW-09						Sampled during baseline event if installed	

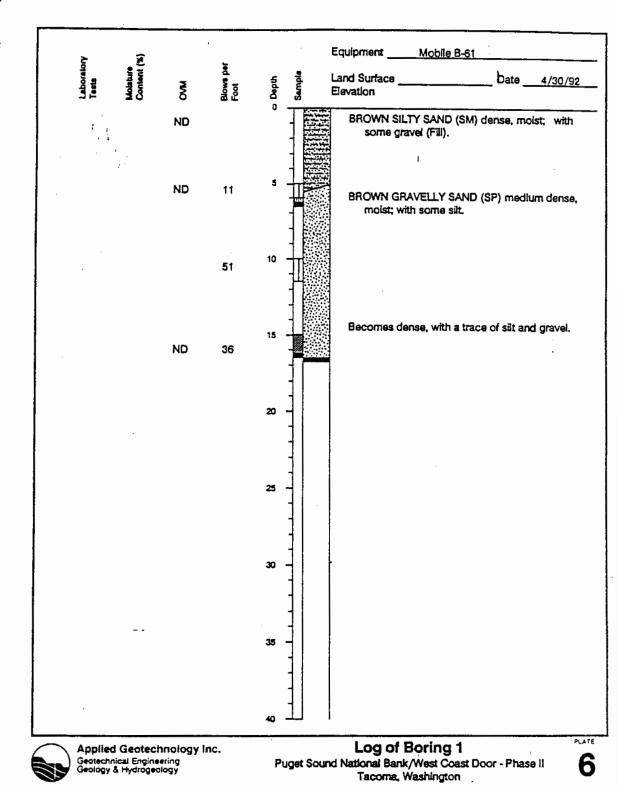
Table 2 Pilot Study Performance Monitoring West Coast Door Site, Tacoma, Washington

		Post Injection Monitoring		ng		
Well ID	Baseline	First	Second	Third	Fourth	Notes
	Sampling	Quarter	Quarter	Quarter	Quarter	
Proposed	Performan	ce Monitor	ing Wells			
MW-18	Х	Х	Х	Х	Х	Water level monitoring during injection event
MW-19	Х	Х	Х	Х	Х	Water level monitoring during injection event
MW-20	Х	Х	Х	Х	Х	Water level monitoring during injection event
MW-21	Х	Х	Х	Х	Х	Water level monitoring during injection event
MW-22	X	Х	Х	Х	Х	Water level monitoring during injection event

Notes:

- 1. X = monitored location.
- 2. Depth to groundwater monitored in all locations during each event and in performance monitoring wells before, during, and after amendment injection.
- 3. All samples analyzed for BTEX and naphthalene by EPA Method 8260D and PAHs by EPA Method 8270E with selected ion monitoring (SIM).
- 4. One trip blank will be analyzed per sample cooler, and one duplicate sample will be analyzed per 20 samples collected.

Soil Boring and Monitoring Well Logs

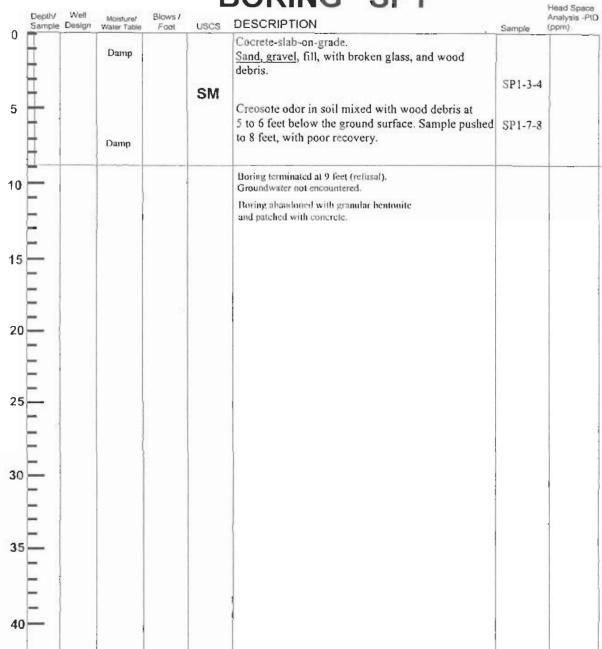


Tacoms, Washington

JOB MUMBER DRAWN APPROVED DATE

15.536.003 SES 18 Jun 92

BORING SP1



Sampler: Continuous Sample collected in 4-foot sections. Driller: ESN-LAR Strateprobe



ENVIRONMENTAL ASSOCIATES, INC.

1380 - 112th Avenue NE, Suite 300 Bellevue, Washington 98004

BORING SP1

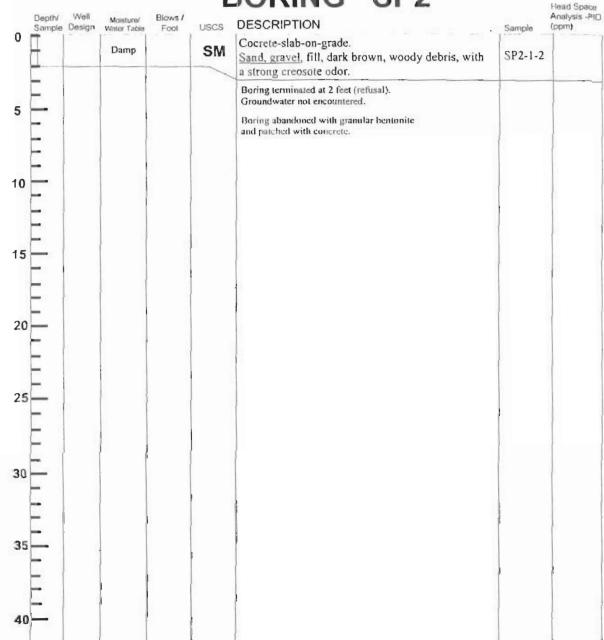
West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number:
JN-26192-1

HC	4				
0	7	12	0	0	6

Logged by:	
RBR	

BORING SP2



Sampler: Continuous Sample collected in 4-loct sections.

Driller: ESN-LAR Strataprobe.



ENVIRONMENTAL ASSOCIATES, INC.

1380 - 112th Avenue NE, Suite 300 Bellevue, Washington 98004

BORING SP2

West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number: JN-26192-1 Date:

07/20/06

Logged by:

Plate:

RBR

A-2

BORING SP3 Head Space Analysis -PID Well Blows / Depth/ Moisture/ Water Table DESCRIPTION Sample Design USCS (ppm) Foot Sample Cocrete-slab-on-grade. Damp Sand, gravel, fill, brick debris, brown. Creosote odor in soil mixed with wood debris at SP3-3-4 3 to 4 feet below the ground surface. SM Sand, gravel, medium to coarse sand, with gravel, 5 no odor below 5 feet. SP3-7-8 Damp SP3-10-11 10 Boring terminated at 11 feet (refusal). Groundwater not encountered. Boring abandoned with granular bentonite and patched with concrete. 20 30 35 40

Sampler: Continuous Sample collected in 4-foot sections.

Driller: ESN-LAR Strataprobe.



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BORING SP3

West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job f	lumber:
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JN-26192-1

Date:

07/20/06

Logged by:

RBR

Plate:

A-3

BORING SP4 Head Space Analysis -PtD (ppm) Well Depth/ Blows / Moisture/ DESCRIPTION Sample Design uscs Water Table Foot Sample Cocrete-slab-on-grade. Sand, gravel, fill, brown. SM Damp Sand, gravel, organic woody debris, dark brown, SP4-3-4 slight creosote odor. Sandy-gravel, medium to coarse sand, mixed with **GM** Damp gravel, brown. No odor. SP4-7-8 Damp SP4-10-11 10 Boring terminated at 11 feet (refusal). Groundwater not encountered. Boring abandoned with granular bentonite and patched with concrete. 20

Sampler: Continuous Sample collected in 4-foot sections. Driller: ESN-LAR Strataprobe.



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BORING SP4

West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number:	Date:	Logged by:	Plate:	Т
JN-26192-1	07/20/06	RBR	A-4	

BORING SP5 Head Space Analysis -PID (ppm) Depth/ Well Sample Design Well Blows / Foot Moisturei Water Table uscs DESCRIPTION Sample Asphalt pavement / gravel base. Damp 0.0 SP5-3-4 **GP** Sandy-Gravel, gravel with medium to coarse sand, and cobbles, brown. SP5-6-7 0.0 Damp SP5-8-9 0,0 Boring terminated at 9 feet (refusal). 10 Groundwater not encountered. Boring abandoned with granular bentonite and patched with asphalt. 20 25 30 35 40

Sampler: Continuous Sample collected in 4-foot sections. Driller: ESN-LAR Strataprobe.



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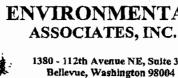
BORING SP5

Former West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number:	Date:	Logged by:	Plate:	
JN-26192-2	09/12/06	RBR	A-1	

BORING SP6 Head Space Analysis -PID Depth/ Well Sample Design Moislure/ Water Table USCS DESCRIPTION Sample Asphalt pavement / gravel base. Damp 0.0 SP6-3-4 GP Sandy-Gravel, gravel with medium to coarse sand, and cobbles, brown. 5 Damp SP6-6-7 0.0 Boring terminated at 7 feet (refusal). Groundwater not encountered. Boring abandoned with granular bentonite 10 and patched with asphalt. 15 20 25 30

Sempler: Continuous Sample collected in 4-foot sections, Drifter: ESN-LAR Strataprobe,



ENVIRONMENTAL ASSOCIATES INC BORING SP6 Former West Coast Door Property

ASSOCIATES, INC.

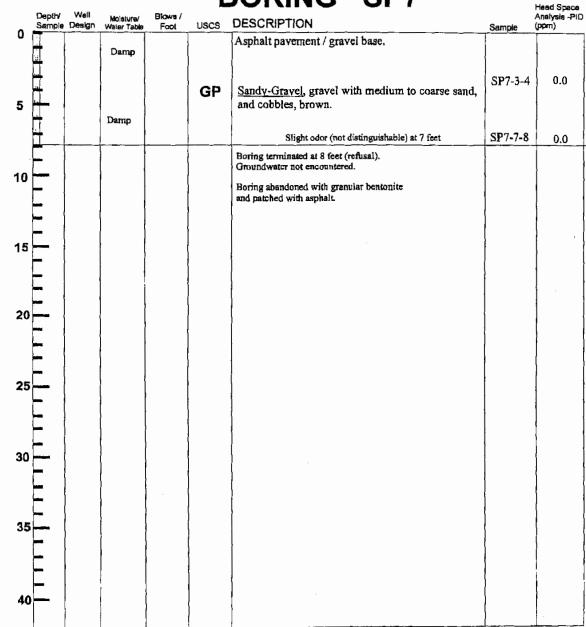
3133 South Cedar Street

1380 - 112th Avenue NE, Suite 300

Tacoma, Washington 98409

		———	
Job Number:	Date:	Logged by:	Plate:
JN-26192-2	09/12/06	RBR	A-2

BORING SP7



Sampler: Continuous Sample collected in 4-foot sections. Driller; ESN-LAR Strataproba,



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BORING SP7

Former West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number:

Date:

Logged by:

JN-26192-2

09/12/06

RBR

A-3

Plate:

BORING SP8 Head Space Analysis -PID Depth/ Well Sample Design Blows / Moisture/ Water Table DESCRIPTION USCS Foot Sample SP Sand, fine to medium sand, with gravel, brown. Damp SP8-3-4 7 70.0 Sandy-Gravel, gravel with medium to coarse sand, 5 GP and cobbles, brown. Damp SP8-7-8 0.0 Boring terminated at 8 feet (refusal). Groundwater not encountered. Boring abandoned with granular bentonite and patched with asphalt. 15 20 25 30

Sampler: Continuous Sample collected in 4-foot sections. Orlifer: ESN-LAR Strateprobe.



35

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BORING SP8

Former West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number:	Date:	Logged by:	Plate:
JN-26192-2	09/12/06	RBR	A-4

BORING SP9 Head Space Analysis -PID (ppm) Depthy Well Sample Design Blows / Foot Moisture/ Water Table DESCRIPTION USCŠ Sample SP Sand, fine to medium sand, organic, with gravel, Damp dark-brown. SP9-3-4 -0.0-Sandy-Gravel, gravel with medium to coarse sand, GP and cobbles, brown. Damp SP9-7-8 0.0 Boring terminated at 8 feet (refusal). Groundwater not encountered. Boring abandoned with granular bentonite and patched with asphalt. 30 35

Sampler: Continuous Sample collected in 4-foot sections. Orliter: ESN-LAR Strateprobe.



ENVIRONMENTAL ASSOCIATES, INC.

1380 - 112th Avenue NE, Suite 300 Bellevue, Washington 98004

BORING SP9

Former West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

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Job Number:	Date:	Logged by:	Plate:
JN-26192-2	09/12/06	RBR	A-5

Date/Time Started : 1-29-07/1040 Date/Time Completed: 1-29-07/1204 LOG OF BORING SB-1 **Total Boring Depth** : 35' Depth to water ATD : 24' (Page 1 of 2) Elevation (ft) : NA **Drilling Method** : HSA Site Name: West Coast Door Property Sampler Type : D+M S.S. 2" diameter 18' : 300 Drive Hammer (lbs) Client: William Swensen PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Recovery **Blow Count** PID (ppm) Samples Graphic Description Sample ID 0 4.0 - 5.5 GRAVEL minor sand, trace silt (80% fine to coarse gravel, 15% fine to coarse sand, 5% silt), brown, moist. 100 20/21/21 6.2 5 08-13-2008 "NPacific-8e185af/public/Project Files/112 Swensen/112-001 West Coast Door Property/Boring Logs/SB1.bo 9.0 - 10.0 GRAVEL minor sand, trace silt (80% fine to coarse gravel, 15% 14/15/19 100 16.2 SB1-9-10 fine to coarse sand, 5% silt), brown-gray, wet. 10 14.0 - 15.5 SAND minor gravel, trace silt (85% fine to coarse sand, 10% fine to coarse gravel, 5% silt), brown, moist. SP 34/30/19 0.0 15 19.0 - 19.5 SAND minor silt, trace gravel (85% fine to coarse sand, 10% silt, 5% coarse gravel), brown, moist. 19.5 - 20.5 Gravelly SAND trace silt (50% medium to coarse sand, 45% SP 100 29-50/5" 70.2 SB1-19-20.5 fine to coarse gravel, 5% silt), brown, moist. SP Drilling Company : Cascade Drilling, Inc. **Drilling Foreman** : James Goble LOG OF BORING SB-1 Equipment : CME Pacific Crest Rep. : Annica Nord (Page 1 of 2)

Date/Time Started : 1-29-07/1040 Date/Time Completed: 1-29-07/1204 LOG OF BORING SB-1 Total Boring Depth : 35' Depth to water ATD : 24' (Page 2 of 2) Elevation (ft) : NA **Drilling Method** : HSA Site Name: West Coast Door Property Sampler Type : D+M S.S. 2" diameter 18' Drive Hammer (lbs) : 300 Client: William Swensen PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Recovery **Blow Count** PID (ppm) Samples Graphic Description Sample ID 20 SP 100 29-50/5" 70.2 SB1-19-20.5 24.0 - 25.5 Gravelly SAND trace silt (50% medium to coarse sand, 45% fine to coarse gravel, 5% silt), brown, wet. SP 90 26-50/6" 2.6 25 08-13-2008 \Pacific-8e185afpublic\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB1.bo 29.0 - 30.5 SAND trace gravel, trace silt (90% fine to coarse sand, 5% fine gravel, 5% silt), brown, wet, assorted cobbles. SP 100 50/6" 34.1 30 34.0 - 35.0 SAND with gravel (80% medium to coarse sand, 20% fine SP 100 22 SB1-34-GW gravel), gray-brown, wet. 35 40 Drilling Company : Cascade Drilling, Inc. Drilling Foreman : James Goble LOG OF BORING SB-1 Equipment : CME Pacific Crest Rep. : Annica Nord (Page 2 of 2)

Date/Time Started : 1-29-07/0818 Date/Time Completed: 1-29-07/1008 LOG OF BORING SB-2 **Total Boring Depth** : 35' Depth to water ATD : 20' (Page 1 of 2) Elevation (ft) : NA **Drilling Method** : HSA Site Name: West Coast Door Property Sampler Type : D+M S.S. 2" diameter 18' Drive Hammer (lbs) : 300 Client: William Swensen PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Recovery **Blow Count** PID (ppm) Samples Description Sample ID 0. 4.0 - 5.5 GRAVEL with sand, trace silt (75% fine to coarse gravel, 20% fine to coarse sand, 5% silt), brown, dry. GP 100 26/28/23 30.1 5 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB2.bo 9.0 - 10.5 GRAVEL with sand, trace silt (75% fine to coarse gravel, 20% fine to coarse sand, 5% silt), brown, dry. GP 10 50/6" 22.3 10 14.0 - 15.5 SAND minor silt (interbedded 100% coarse sand with 90% coarse sand, 10% silt), brown, moist. SP 100 29-50/6" 24.6 SB2-14-15.5 15 19.0 - 20.5 SAND minor silt, trace gravel (85% medium to coarse sand, 10% silt, 5% fine to coarse gravel), brown, wet. SP 100 36-50/6" 10.6 SB2-19-20.5 Drilling Company : Cascade Drilling, Inc. Drilling Foreman : James Goble LOG OF BORING SB-2 Equipment : CME Pacific Crest Rep. : Annica Nord (Page 1 of 2)

Date/Time Started : 1-29-07/0818 Date/Time Completed: 1-29-07/1008 LOG OF BORING SB-2 Total Boring Depth : 35' Depth to water ATD : 20' (Page 2 of 2) Elevation (ft) : NA **Drilling Method** : HSA Site Name: West Coast Door Property Sampler Type : D+M S.S. 2" diameter 18' Drive Hammer (lbs) : 300 Client: William Swensen PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Recovery **Blow Count** PID (ppm) Samples Description Sample ID % 20 SP 36-50/6" 10.6 SB2-19-20.5 24.0 - 25.0 SAND (100% fine to medium sand), brown, wet. SP 100 35-50/6" 29.7 25 08-13-2008 \Pacific-8e185afpublic\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB2.bo 29.0 - 30.5 SAND trace gravel, trace silt (90% fine to coarse sand, 5% fine to coarse gravel, 5% silt), gray, wet. SP 100 26-50/6" 107 30 34.0 - 35.0 GRAVEL minor sand, trace silt (85% fine to coarse 100 15.1 SB2-34-GW gravel, 10% fine to coarse sand, 5% silt), gray, wet. 35 Drilling Company : Cascade Drilling, Inc. Drilling Foreman : James Goble LOG OF BORING SB-2 Equipment : CME Pacific Crest Rep. : Annica Nord (Page 2 of 2)

LOG OF BORING SB-3

(Page 1 of 2)

Site Name: West Coast Door Property

Client: William Swensen

Project #: 112 001

Date/Time Started : 1-31-07/1141 Date/Time Completed: 1-31-07/1416

Total Boring Depth : 33' Depth to water ATD : 25' Elevation (ft) : NA **Drilling Method** : HSA

Sampler Type : D+M S.S. 2" diameter 18'

Drive Hammer (lbs) : 140

15 B	ENDI	BEND. WA 98045	Project #: 112-001						
Depth In Feet	Samples	Desc	ription	nscs	Graphic	% Recovery	Blow Count	PID (ppm)	Sample ID
0—		0.0 - 5.0 Soil cuttings are rounded or	obbles with some sand and silt.	GP					
10 — - - - - -	X	10.0 - 11.5 SAND with silt, minor gra 20% silt, 10% coarse gravel), brown	vel (70% fine to medium sand, dry, no odor.	SP		5	50/1"	49.2	SB3-10-11.5
	X	15.0 - 16.5 GRAVEL minor sand, tra 10% fine to coarse sand, 5% silt), br	ce silt (85% fine to coarse gravel, own, moist, no odor.	GP		40	50/3"	35.7	SB3-15-16.5
Drilling									

Drilling Foreman

08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB3.bo

Equipment : CME 65 Pacific Crest Rep. : Annica Nord

: Steve Choate

LOG OF BORING SB-3

(Page 1 of 2)

LOG OF BORING SB-3

(Page 2 of 2)

PACIFIC CREST ENVIRONMENTAL 1933 Bendigo Boulevaro North PD 80x 992 N 0 8 T 0 4 2 9 8 0 4 2 9

08-13-2008 NPacific-8e185aflpublic/Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB3.bo

Pacific Crest Rep.

: Annica Nord

Site Name: West Coast Door Property

Client: William Swensen

Date/Time Started : 1-31-07/1141 Date/Time Completed : 1-31-07/1416

Total Boring Depth : 33'
Depth to water ATD : 25'
Elevation (ft) : NA
Drilling Method : HSA

Sampler Type : D+M S.S. 2" diameter 18'

(Page 2 of 2)

Drive Hammer (lbs) : 140

PACI		DREST ENVIRONMENTAL D BOULEVARD NORTH PO BOX 952 B E N D , W A 9 B D 4 5	Project #: 112-001						
Depth In Feet	Samples	Desc	nscs	Graphic	% Recovery	Blow Count	PID (ppm)	Sample ID	
20-	:	20.0 - 21.5 NO RECOVERY.				0	50/1"	_	-
		25.0 - 26.5 NO RECOVERY. Samples of the second seco		-		0	50/0"	-	-
30-		30.0 - 31.5 NO RECOVERY.				0	50/1"	-	-
-		33.0 - 34.5 GRAVEL with minor sand gravel, 10% fine to coarse sand, 10%	d, minor silt (80% fine to coarse 6 silt), brown, wet, no odor.	GP		5	50/3"	96.6	SB3-33-GW
35—									
Drilling Drilling Equipm	Foren					L	OG OF	BOR	ING SB-3

LOG OF BORING SB-4

(Page 1 of 2)

Site Name: West Coast Door Property

Client: William Swensen

Date/Time Started : 5-11-07/0949 Date/Time Completed: 5-11-07/1135

Total Boring Depth : 35' Depth to water ATD : 25' Elevation (ft) : NA **Drilling Method** : HSA

Sampler Type : D+M S.S. 2" diameter 18'

Drive Hammer (lbs) : 300

PACIFIC CREST ENVIRONMENTAL

1533 BENDIGO BOULEVARD NORTH PO BOX 952

Project #: 112-001 Depth In Feet Recovery **Blow Count** PID (ppm) Graphic Description Sample ID 0 5 5.0 - 6.5 Sandy GRAVEL minor silt (65% fine to coarse gravel, 25% fine to coarse sand, 10% silt), brown, moist, no odor. GΡ 60 50/6" 16.5 10 10.0 - 10.5 Silty SAND (70% fine to coarse sand, 30% silt), SM brown, moist, no odor. 100 20/25/22 SB4-10-11.5 16.2 SP 10.5 - 11.5 SAND trace silt (95% fine to medium sand, 5% silt), brown, moist, no odor. 15 15.0 - 16.5 SAND with gravel with silt (70% fine to coarse sand, 20% fine to coarse gravel, 10% silt), brown, moist, no SP 41-50/6" 15.5 odor. 20

Drilling Company Drilling Foreman

08-13-2008 NPacific-8e185afipublic\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB4.bo

: Cascade Drilling, Inc.

Equipment : CME Pacific Crest Rep. : Annica Nord

: James Goble

LOG OF BORING SB-4

(Page 1 of 2)

Date/Time Started : 5-11-07/0949 Date/Time Completed: 5-11-07/1135 LOG OF BORING SB-4 **Total Boring Depth** : 35' Depth to water ATD : 25' (Page 2 of 2) Elevation (ft) : NA **Drilling Method** : HSA Site Name: West Coast Door Property Sampler Type : D+M S.S. 2" diameter 18' : 300 Drive Hammer (lbs) Client: William Swensen PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Recovery **Blow Count** PID (ppm) Graphic Description Sample ID 20 20.0 - 21.5 SAND with gravel with silt (70% fine to coarse sand, 20% fine to coarse gravel, 10% silt), brown, moist, no SP 35 50/6" 30.6 SB4-21-21.5 odor. Cobble at bottom of sampler. 25 25.0 - 25.75 SAND with gravel with silt (70% fine to coarse SP sand, 20% fine to coarse gravel, 10% silt), brown, moist, no 40 50/6" 27.0 SP 25.75 - 26.5 SAND trace silt (95% fine sand, 5% silt), brown, wet, no odor. 08-13-2008 NPacific-8e185afipublic\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB4.bo 30 30.0 - 31.5 SAND trace silt (95% fine to medium sand, 5% silt), brown, wet, no odor. Cobble at the bottom of the sampler. SP 50/6" 35 25.1 35 35.0 - 36.5 SAND (100% fine to medium sand), brown, wet, no odor. SP 50/6" 21.2 SB4-35-GW Drilling Company : Cascade Drilling, Inc. **Drilling Foreman** : James Goble LOG OF BORING SB-4 Equipment : CME Pacific Crest Rep. : Annica Nord (Page 2 of 2)

Date/Time Started : 5-11-07/1301 Date/Time Completed: 5-11-07/1425 LOG OF BORING SB-5 **Total Boring Depth** : 35' Depth to water ATD : 20' (Page 1 of 2) Elevation (ft) : NA **Drilling Method** : HSA Site Name: West Coast Door Property Sampler Type : D+M S.S. 2" diameter 18' Drive Hammer (lbs) Client: William Swensen : 300 PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Recovery **Blow Count** PID (ppm) Graphic Description Sample ID 0 5 5.0 - 6.5 Sandy GRAVEL minor silt (70% fine to coarse gravel, 20% fine to coarse sand, 10% silt), brown, moist, no odor. GΡ 17/18/21 34.5 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB5.bo 10 10.0 - 10.5 Sandy GRAVEL minor silt (70% fine to coarse SM gravel, 20% fine to coarse sand, 10% silt), brown, moist, no 40-50/6" 55 35.8 SM odor. Large cobble. 10.5 - 11.5 Silty SAND (70% fine sand, 30% silt), brown, moist, no odor. 15 15.0 - 15.5 Silty SAND minor gravel (60% fine to coarse sand, 30% silt, 10% fine gravel), brown, moist, no odor. SM 46-50/2" 39.6 SB5-15-16.5 SP 15.5 - 16.5 SAND with gravel, trace silt (70% fine to coarse sand, 25% fine to coarse gravel, 5% silt), brown, moist, no odor. 20 **Drilling Company** : Cascade Drilling, Inc. Drilling Foreman : James Goble LOG OF BORING SB-5 Equipment : CME

(Page 1 of 2)

Pacific Crest Rep.

: Annica Nord

Date/Time Started : 5-11-07/1301 Date/Time Completed: 5-11-07/1425 LOG OF BORING SB-5 **Total Boring Depth** : 35' Depth to water ATD : 20' (Page 2 of 2) Elevation (ft) : NA **Drilling Method** : HSA Site Name: West Coast Door Property Sampler Type : D+M S.S. 2" diameter 18' Drive Hammer (lbs) Client: William Swensen : 300 PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Recovery Blow Count PID (ppm) Graphic Description Sample ID 20 20.0 - 21.5 Gravelly SAND minor silt (45% fine to coarse sand, 45% fine to coarse gravel, 10% silt), brown, wet, no odor. SP 100 40-50/6" 36.6 25 25.0 - 25.5 Gravelly SAND minor silt (45% fine to coarse sand, SP 45% fine to coarse gravel, 10% silt), brown, wet, no odor. 29-50/6" 100 32.6 SM 25.5 - 26.5 Silty SAND trace silt (65% fine sand, 35% silt), brown, wet, no odor. 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB5.bo 30 30.0 - 30.5 Gravelly SAND minor silt (45% fine to coarse sand, SP 45% fine to coarse gravel, 10% silt), brown, wet, no odor. GP 50/6" 60 42.2 30.5 - 31.0 GRAVEL minor sand, minor silt (70% fine gravel, 15% fine to coarse sand, 15% silt), brown, wet, no odor. 35 35.0 - 36.5 Gravelly SAND trace silt (60% fine to coarse sand, 35% fine gravel, 5% silt), gray, wet, creosote odor. SP 50/6" 43.8 SB5-35-36.5 SB5-35-GW

Drilling Company **Drilling Foreman**

Pacific Crest Rep.

: Cascade Drilling, Inc. : James Goble

Equipment : CME

: Annica Nord

LOG OF BORING SB-5

(Page 2 of 2)

Date/Time Started : 4-01-2009 / 8:18 Date/Time Completed : 4-01-2009 / 4:28 LOG OF BORING SB-6 Total Boring Depth : 39 feet Depth to water ATD : 20 feet (Page 1 of 1) Elevation (ft) : NA **Drilling Method** : Direct Push & HSA Site Name: West Coast Door Sampler Type : Macro-Core & piston sampler Client: Swensen Enterprises, LLC Project #: 112-001 Depth In Feet % Recovery PID (ppm) Graphic Description Sample ID No Sampling above 28 feet bgs. From auger cuttings: From Approximately 0 - 12: GRAVEL and coarse to medium sand GP 10 From Auger Cuttings: From Approximately 12 -28: SAND trace silt, trace gravel (90% fine to medium grained sand, 5% silt, 5% gravel), medium brown, moist, no odor. \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB6.bor SP 20 SB6-22-RGW 28 -31.5 SAND (100% medium to coarse grained sand), light brown to medium gray, wet, no odor. SP 30 100 0.0 31.5 - 32 SAND minor silt (85% fine sand, 15% silt), medium gray, wet, no 32 - 32.5 SAND (100% fine sand) medium gray, wet, no odor. SB6-37-RGW 32.5 - 35 Silty SAND (65% fine sand, 35% silt), medium gray, moist, no odor. 100 0.1 SM 35 - 39 Sandy SILT (60% SILT, 40% fine sand), medium gray, moist, no odor. ML 100 0.3 Bottom of Boring at 39 feet bgs **Drilling Company** : ESN Northwest **Drilling Foreman** : Noel Knopf LOG OF BORING SB-6 Equipment : AMS Powerprobe 9630 Pacific Crest Rep. : Monty Busbee

(Page 1 of 1)

Date/Time Started : 4-01-2009 / 8:45 Date/Time Completed : 4-02-2009 / 4:28 LOG OF BORING SB-9 Total Boring Depth : 43 feet Depth to water ATG : 24.13 (Page 1 of 2) Elevation (ft) : NA **Drilling Method** : Direct Push Site Name: West Coast Door Sampler Type : screen-point groundwater Client: Swensen Enterprises, LLC : Macro-core Project #: 112-001 **Depth In Feet** Recovery PID (ppm) Graphic Description Sample ID 0 - 4 Sandy GRAVEL with silt (50% gravel, 30% fine to coarse sand, 20% silt), light to dark gray, dry, no odor GP-GM 70 0.0 4 - 8 As above GP-GM 80 0.0 8 - 12 As above GP-GN 10 50 0.0 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB9.bo 12 - 12.5 SAND (100% coarse sand), medium brown, moist, no odor. 12.5 - 16 Sandy GRAVEL with silt (50% gravel, 30% fine to coarse sand, 20% silt), light to dark gray, moist, no odor. 60 0.0 GP-GN 16 - 19 As above. GP-GM 100 0.0 19 - 20 SAND to Sandy Silty GRAVEL (Slough?) medium gray, moist, no odor. 3P-SM 20 - 21 GRAVEL with sand to SAND (Slough *4 foot sampler filled from one 20 foot of advancement). GP 100* 0.0 21 - 22 As above. 100* 0.0 Broken probe rods and sampler. Advancing boring with hollow stem augers to 40 feet. SB9-32-RGW

Drilling Company : ESN Northwest **Drilling Foreman** : Noel Knopf

Equipment : AMS Powerprobe 9630 Pacific Crest Rep. : Monty Busbee

LOG OF BORING SB-9

(Page 1 of 2)

Date/Time Started : 4-01-2009 / 8:45 LOG OF BORING SB-9 Date/Time Completed : 4-02-2009 / 4:28 Total Boring Depth : 43 feet Depth to water ATG : 24.13 (Page 2 of 2) Elevation (ft) : NA Drilling Method : Direct Push Site Name: West Coast Door Sampler Type : screen-point groundwater Client: Swensen Enterprises, LLC : Macro-core Project #: 112-001 Depth In Feet % Recovery PID (ppm) Samples Graphic Description Sample ID 30 SB9-32-RGW 40 - 41 SAND (100% fine to coarse sand), medium gray, wet, VOC odor. 40 41 - 42 GRAVEL with sand (70% gravel, 30% fine to medium sand), medium SP gray, wet, VOC odor. 100 4.6 GP \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\SB9.bor 42 - 43 Silty SAND (60% fine to medium sand, 40% silt), medium gray, wet, SM slight VOC odor. GM 100 SB9-43-RGW 3.1 43 - 43.5 GRAVEL trace silt (95% gravel, 5% silt), medium gray, moist, VOC SM 43.5 - 45 Silty SAND (60% fine to medium sand, 40% silt), medium gray, SP moist, VOC odor. 45 - 45.5 SAND trace silt (95% fine to medium sand, 5% silt), medium gray, 100 moist, VOC odor. SM 45.5 - 49 Silty SAND (60% fine to medium sand, 40% silt), medium gray, moist, VOC odor. Bottom of Boring at 49 feet bgs 50 **Drilling Company** : ESN Northwest **Drilling Foreman** : Noel Knopf LOG OF BORING SB-9 Equipment : AMS Powerprobe 9630 Pacific Crest Rep. : Monty Busbee (Page 2 of 2)



Coordinate System: NAD83/98

Ground Surf Elev. and Datum245.9 NGVD 29

Latitude/Northing: 697992.86 Longitude/Easting: 1149534.32

Boring Location:

Drill Date: June 11, 2010 **Logged By:** John LaManna

Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA **Client:** Bill Swensen **Sample Method:** 18" D&M Split-spoon **Project:** Swensen-WCD

Boring ID: SP2-B

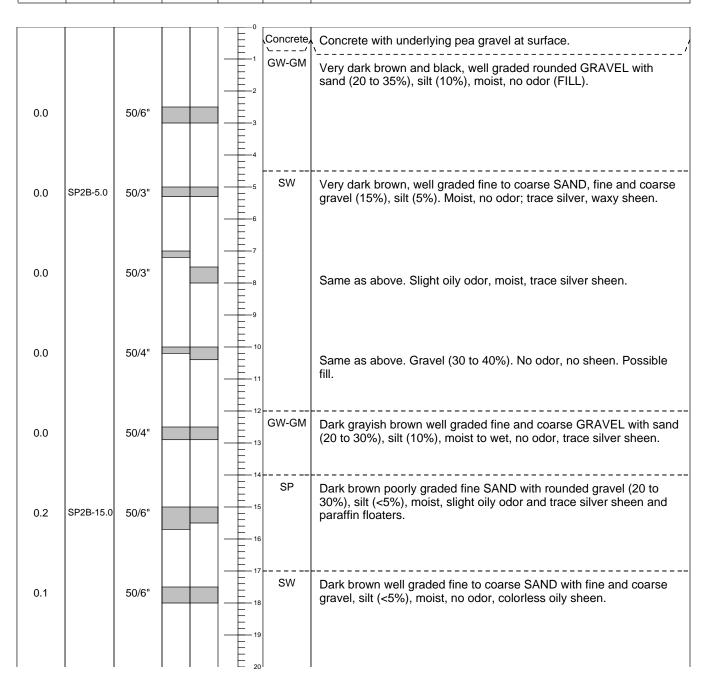
Boring Diameter: 8" Task:

Boring Depth (ft bgs): 46 ft bgs Address: 3133 Cedar St,

Groundwater ATD (ft bgs): 24 Tacoma, WA.

Remarks: Boring backfilled with bentonite chips to 5 feet bgs, then concrete patched.

L							
	PID	SAMPLE	BLOW	DRIVEN /	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS
	(ppm)	ID	COUNT	RECOVERED	FT BGS	SYMBOL	





Coordinate System: NAD83/98

Ground Surf Elev. and Datum245.9 NGVD 29 Sample Method: 18" D&M Split-spoon

Latitude/Northing: 697992.86 Longitude/Easting: 1149534.32

Boring Location:

Drill Date: June 11, 2010 Logged By: John LaManna

Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA Client: Bill Swensen
Sample Method: 18" D&M Split-spoon Project: Swensen-WCD

Boring ID: SP2-B

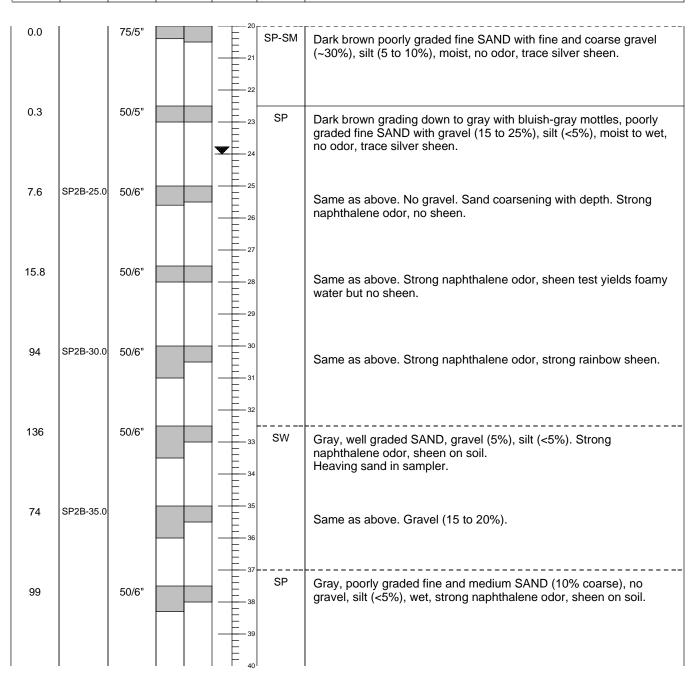
Boring Diameter: 8" Task:

Boring Depth (ft bgs): 46 ft bgs Address: 3133 Cedar St,

Groundwater ATD (ft bgs): 24 Tacoma, WA.

Remarks: Boring backfilled with bentonite chips to 5 feet bgs, then concrete patched.

PID	SAMPLE	BLOW	DRIVEN /	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS
(ppm)	ID	COUNT	RECOVERED	FT BGS	SYMBOL	



= denotes groundwater table



Coordinate System: NAD83/98

Ground Surf Elev. and Datum245.9 NGVD 29 Sample Method: 18" D&M Split-spoon

Latitude/Northing: 697992.86

Longitude/Easting: 1149534.32

Boring Location:

Drill Date: June 11, 2010 **Logged By:** John LaManna

Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA

Client: Bill Swensen

Sample Method: 18" D&M Split-spoon

Project: Swensen-WCD

Boring ID: SP2-B

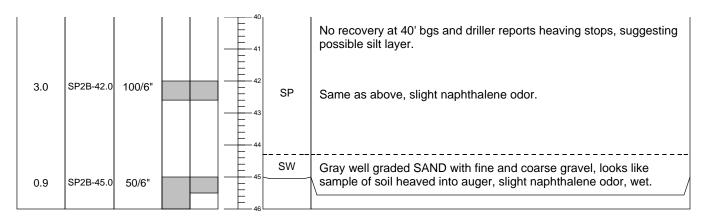
Boring Diameter: 8" Ta

Boring Depth (ft bgs): 46 ft bgs Address: 3133 Cedar St,

Groundwater ATD (ft bgs): 24 Tacoma, WA.

Remarks: Boring backfilled with bentonite chips to 5 feet bgs, then concrete patched.

PID	SAMPLE	BLOW	DRIVEN /	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS
(ppm)	ID	COUNT	RECOVERED	FT BGS	SYMBOL	



FLOVELCNIER	PROJECT: West Coast Door	LOCATION: 31		BORING	D: SB-40		
FLOYDISNIDER	LOGGED BY:		St, Tacoma WA BORING LOCATION:				
strategy • science • engineering	K. Anderson		warehouse no	orth of form	ner retort		
DRILLED BY:	IX. / IIIdel 3011	NORTHING:		ASTING:	TICL TOTAL		
Zack Bailey, Holocene Drilling		698011.38		1149553.	06		
DRILLING EQUIPMENT: Sonic LAR		SURFACE ELEVATION: 2	E 2 1	COORDINATE SYSTEM: NAVD88			
DRILLING METHOD:		TOTAL DEPTH (ft bgs):		/ATER (ft bgs):		
5' x 4" Rods		75		34			
SAMPLING METHOD/SAMPLER LENGTH: 4" Liner Bag		BORING DIAME 4" inner/6" ou		10/1/2018			
(feet) Symbol (color, texture, moisture, MAJC	escription and Observations OR CONSTITUENT, odor, staining	g, sheen, debris, etc.)	Drive/ Recovery	PID (ppm)	Sample ID		
6" of CONCRETE .							
Gray-brown, well-graded S A no odor. At 2 feet bgs, becomes brown At 2.5 feet bgs, becomes date at 2.5 feet bgs, becomes light At 5 feet bgs, cobbles up	vn. ark brown gray-brown and dry.			0.2			
Well-graded GRAVEL with Poorly-graded SAND with g	,	/ sand, moist.		0.0			
Brown, well-graded SAND of the second seco				0.0			
ABBREVIATIONS:		NOTES:		<u> </u>			
ft bgs = feet below ground surface USCS = Unified	Soil Classification System s groundwater table	Recovered intervals evenly	y decompressed fo	or log unless	otherwise noted.		

		PROJECT:	LOCATION:	3102 S Cedar	BORING II	D: CD 40	
FL	OYDISNIDER	West Coast Door		St, Tacoma WA		SB-40	
	egy • science • engineering	LOGGED BY:		BORING LOCATION:			
	5 57	K. Anderson			north of former retort		
DRILLEI Zack	ви: Bailey, Holocene Drilling		NORTHING : 698011.38		E ASTING : 1149553.0)6	
	IG EQUIPMENT:		SUDFACE	C	OORDINATE		
Sonic			ELEVATION:	252.1	NAVD88	201012	
DRILLIN	IG METHOD:		TOTAL DEPTH	H (ft bgs): D	EPTH TO W	ATER (ft bgs):	
5' x 4'	' Rods		75		34		
	ING METHOD/SAMPLER LENGTH:		BORING DIAM 4" inner/6"		DRILL DATE : 10/1/2018		
	er Bag		4 111116170				
Depth (feet)		scription and Observations R CONSTITUENT, odor, staining	, sheen, debris, etc.)		PID (ppm)	Sample ID	
20	Silty SAND with fine gravel, some standard sand with gravel. Well-graded SAND with gravel.		no odor.		0.0		
28 —	Gray-brown poorly-graded m to wet, no odor. At 27.5 feet bgs, slight creos		e fine gravel, moist	-	0.0		
32 —	From 29.5-30 feet bgs, slight brown film in sheen test. At 30 feet bgs, moderate cre and brown film in sheen test.	osote odor present. Lig				SB-40-29.5-30 SB-40-31.5-32	
-	Well-graded SAND with large odor and slight brown film in SW: At 35 feet, no odor and no sh	sheen test.	et. Slight creosote		1.2		
36 —	Gray-brown fine silty SAND , Gray-brown poorly-graded fine		· ·		0.7		
	SP: VIATIONS: = feet below ground surface. USCS = Unified.		NOTES: Recovered intervals eve	enly decompressed for		SB-40-39-39.5	
	= feet below ground surface USCS = Unified = parts per million = denotes	Soil Classification System groundwater table	Mecovered littervals eve	any decompressed it	n log utiless (ou ici wise Hüleü.	

	0.1/	D I CALLED ED	PROJECT:	LOCATION: 31		BORI	NG ID: SB-40
F L	OY	DISNIDER	West Coast Door		, Tacoma WA		3D-40
strat	egy •	science • engineering	LOGGED BY: K. Anderson	BORING LOCAT		orth of	formar ratort
DRILLEI	n pv.		K. Affuersoff	NORTHING:	warehouse no	ASTINO	
		Holocene Drilling		698011.38	-	11495	
DRILLIN	IG EQUIPI	MENT:		SURFACE	С	OORDII	NATE SYSTEM:
Sonic	LAR			ELEVATION: 2	252.1	NAVD	88
DRILLIN	IG METHO	DD:		TOTAL DEPTH ((ft bgs):		O WATER (ft bgs):
	' Rods			75		34	
	I NG MET H er Bag	IOD/SAMPLER LENGTH:		BORING DIAME 4" inner/6" or	I .	10/1/2	
Depth	USCS	Sail Day	scription and Observations	4 11110170 00	Drive/	PID	
(feet)	Symbol	(color, texture, moisture, MAJO		g, sheen, debris, etc.)		(ppm)	Sample ID
40							
		Well-graded SAND with silt a	and gravel no odor or	sheen		0.0	
			and graver, no odor or	SHOOM.			
_	SW-SM						
						0.0	
		Dark gray poorly-graded med trace silt, no odor or sheen.	dium SAND with fine r	ounded gravel and		0.0	
44 —		trace siit, no odor or sneen.					
_	SP						
						0.0	
48 —		Gray, very firm SILT, slightly	moist.				
	: : : SM :	Well-graded silty SAND , no	odor.			0.0	
_		Well-graded GRAVEL with to	raco sand and silt no	odor			
	(GW)	weii-graded GRAVEL with the	face Sand and Sill, no	odoi.			
	0 • 0						
52 —		Brown, well-graded SAND w	ith silt and gravel, moi	st to wet, no odor or			
32 —		sheen.				0.0	
	C)A/ C)A						
	SW-SM						
_							
	0 • 0	Brown, well-graded GRAVEI	L with trace sand and	silt, wet, no odor.			
56 —) · · · ·						
	(GW)						
	0.0						
		Brown, well-graded SAND w	rith gravel and trace sil	lt, no odor.		00	
_]::::::::					0.0	
60	VIATIONS			NOTES:			
ft bgs	VIATIONS = feet belo	ow ground surface USCS = Unified		Recovered intervals evenl	y decompressed for	or log un	lless otherwise noted.
ppm =	parts per	million = denotes	groundwater table				

t Coast Door	N Tachma WA		NG ID: SB-40
ED BY: BORING LO	St, Tacoma WA		
	perty warehouse n	orth of	former retort
NORTHING	6 :	EASTING	G:
698011.	38	11495	53.06
SURFACE ELEVATION	N: 252.1	COORDII NAVD	NATE SYSTEM: 88
TOTAL DEI	PTH (ft bgs):		O WATER (ft bgs):
75		34	
	1	10/1/2	
and Observations ITUENT, odor, staining, sheen, debris, etc.)	Drive/ Recovery	PID (ppm)	Sample ID
and gravel, wet, no odor.		0.0	
gray-brown, firm sandy SILT			SB-40-68-69
t a	698011. SURFACE ELEVATIO TOTAL DE 75 BORING D	SURFACE ELEVATION: 252.1 TOTAL DEPTH (ft bgs): 75 BORING DIAMETER: 4" inner/6" outer n and Observations STITUENT, odor, staining, sheen, debris, etc.) Torve/Recovery and gravel, wet, no odor.	SURFACE ELEVATION: 252.1 COORDII NAVD TOTAL DEPTH (ft bgs): 75 34 BORING DIAMETER: 4" inner/6" outer 10/1/2 n and Observations STITUENT, odor, staining, sheen, debris, etc.) PID (ppm) t and gravel, wet, no odor.

FLOYDISNIDER	PROJECT: West Coast Door	LOCATION: 31	02 S Cedar Tacoma WA	BORIN	SB-41	
strategy • science • engineering	LOGGED BY:	BORING LOCAT				
strategy • science • engineering	K. Anderson	Site property warehouse east of former retort				
DRILLED BY:		NORTHING:		ASTING		
Zack Bailey, Holocene Drilling		697981.38		114958		
DRILLING EQUIPMENT: Sonic LAR		SURFACE ELEVATION: 2	E2 1	NAVD8	IATE SYSTEM: 38	
DRILLING METHOD:		TOTAL DEPTH (• .		O WATER (ft bgs):	
5' x 4" Rods		85		31		
SAMPLING METHOD/SAMPLER LENGTH: 4" Liner Bag		BORING DIAMET 4" inner/6" ou		10/2/20		
Depth USCS (feet) Symbol (color, texture, moisture, MAJO	scription and Observations R CONSTITUENT, odor, stainin	g, sheen, debris, etc.)		PID (ppm)	Sample ID	
0 6" of CONCRETE .						
OGP: Rounded GRAVEL.	ND with amount rounder	d arough and little ailt				
Gray-brown, well-graded SA dry to moist, no odor.	with Small rounded	a gravei and little siit,				
Black-brown, well-graded S . creosote odor.	AND with silt and woo	d fragments. Slight				
				0.0		
					SB-41-4-5	
Λt 5 feet has no wood fragm						
At 5 feet bgs, no wood fragments or odor present.						
SW-SM						
				0.0		
Brown, poorly-graded mediu	m SAND , slight musty	odor. At 10.5 feet		0.0		
bgs, becomes light brown.				0.0		
SP At 11.5 feet bgs, 2" lense of	dark brown sand.	-	-			
12 — 12 — 12 12 12 12 12 12 12 12 12 12 12 12 12						
Brown, well-graded SAND w	ith silt and gravel, moi	st, no odor or sheen.		0.0		
16 — (1) (1)				0.0		
[[2522]						
				0.0		
ABBREVIATIONS:		NOTES:				
ft bgs = feet below ground surface USCS = Unified	Soil Classification System groundwater table	Recovered intervals evenly	y decompressed fo	or log unl	ess otherwise noted.	

FLOYD	SNIDER	PROJECT: West Coast Door	LOCATION: 3	102 S Cedar t, Tacoma WA	BORING	SB-41
strategy • science		LOGGED BY:	BORING LOCA			
strategy . science	e • engineering	K. Anderson	Site propert	y warehouse ea	st of for	mer retort
DRILLED BY:			NORTHING:	E	ASTING:	
Zack Bailey, Holocene	e Drilling		697981.38	697981.38 1149583.06		
DRILLING EQUIPMENT:			SURFACE	CEO 1	OORDINA	ATE SYSTEM:
Sonic LAR			ELEVATION:	252.1	NAVD8	8
DRILLING METHOD:			TOTAL DEPTH			WATER (ft bgs):
5' x 4" Rods			85		31	
SAMPLING METHOD/SAMF 4" Liner Bag	PLER LENGTH:		BORING DIAMETER 4" inner/6" outer		RILL DAT 10/2/20	
Depth USCS	Call Day	and the second Observations	4 11110170 0			
	solor, texture, moisture, MAJOI	scription and Observations R CONSTITUENT, odor, staining, she	en, debris, etc.)	_	PID (ppm)	Sample ID
depth. I At 24.5 light rai At 25 fe	No sheen. feet bgs, moderate o nbow sheen in sheen eet bgs, trace brown o	ray with slight odor. Odor in dor present and brown oily test. bily film and sheen present and present and present and sheen present and sheen present and sheen or o	droplets and in sheen test.		1.1 0.6 — 2.3 —	SB-41-24-25 SB-41-25.5-26 SB-41-28.5-29
At 31 fe 32 — SM Gray si Gray, w SW-SM Gray, p SP At 38 fe	lty SAND , wet, no odd vell-graded SAND with oorly-graded fine SA I	ND, wet. Slight crosote odd	or or sheen. or on water. No .		0.0	
ABBREVIATIONS:	surface LISCS - Unified	NOT	ES: covered intervals even	ly decompressed fo	or log uplo	se atherwise noted

	OVDICNIDED	PROJECT: West Coast Door	LOCATION: 3102	2 S Cedar acoma WA	BORING	SB-41
	OYD SNIDER egy • science • engineering	LOGGED BY:	BORING LOCATIO			
Strati	egy • science • engineering	K. Anderson	Site property w		st of form	er retort
DRILLEI Zack I	DBY: Bailey, Holocene Drilling		NORTHING: 697981.38		ASTING: 1149583.	06
	G EQUIPMENT:		SURFACE			E SYSTEM:
Sonic LAR ELI			ELEVATION: 252	7 1	NAVD88	E STSTEM.
	G METHOD:		TOTAL DEPTH (ft I	•	EPTH TO V 31	/ATER (ft bgs):
5' x 4" Rods SAMPLING METHOD/SAMPLER LENGTH:			BORING DIAMETE		RILL DATE	<u>:</u>
4" Line			4" inner/6" oute		10/2/2018	3
Depth (feet)		scription and Observations R CONSTITUENT, odor, staining, sheen,	debris, etc.)		PID (ppm)	Sample ID
40	SP-SM Gray-brown, well-graded SA sheen.	ND with silt and gravel, wet,	no odor or		0.0	
48 —	Gray-brown, well-graded SA	ND with little silt, no odor or s	sheen.		0.0	
_	Gray, very firm SILT , slightly	moist.	oist, no odor.		0.0	SB-41-49-50
52 —	GW, 6				0.0	
_		velly SAND , moist. ndy GRAVEL with silt, moist,	no odor.		0.0	
56 —	Gray-brown, well-graded gra	velly SAND with silt, wet, no	odor.		0.0	
	/IATIONS:	NOTES				
ft bgs ppm =	= feet below ground surface USCS = Unified parts per million = denotes	Soil Classification System Recovers groundwater table	rered intervals evenly o	lecompressed fo	or log unless	otherwise noted.

FLOVDICNIDED	PROJECT: West Coast Door		3102 S Cedar	BORI	SB-41
FLOYDISNIDER	LOGGED BY:	BORING LOCA	St, Tacoma WA		3D-41
strategy • science • engineering	K. Anderson		ty warehouse e	ast of f	ormer retort
DRILLED BY:		NORTHING:			G:
Zack Bailey, Holocene Drilling		697981.38		11495	
DRILLING EQUIPMENT: Sonic LAR		SURFACE ELEVATION:	252.1	COORDINATE SYSTEM: NAVD88	
DRILLING METHOD:		TOTAL DEPTH	I (ft bgs):	DEPTH TO WATER (ft bgs):	
5' x 4" Rods		85	ETED.	31	
SAMPLING METHOD/SAMPLER LENGTH: 4" Liner Bag		4" inner/6"		DRILL DATE: 10/2/2018	
Depth USCS Soil Detection (feet) Symbol (color, texture, moisture, MAJO)	scription and Observations R CONSTITUENT, odor, staining, sheen, o	debris, etc.)	Drive/ Recovery	PID (ppm)	Sample ID
GW. Gray-brown, well-graded GR Gray-brown, well-graded gra -SW-SM Gray-brown, well-graded silty 64 — SM	velly SAND with silt, wet, no o	odor.		0.0	
Gray-brown, well-graded gra	velly SAND with silt, wet, no o	odor.		0.0	
Gray-brown, poorly-graded r	nedium SAND , no odor.			0.0	
76 — SM SM SO STATE OF THE STAT	NOTES			0.0	
ABBREVIATIONS:	NOTES:	orad intarvala ava	nly docomprossed t	for loa ur	alaca athanuica natad

	YD SNIDER - science - engineering		St	02 S Cedar , Tacoma W	Α	SB-41	
strategy		LOGGED BY:	BORING LOCAT	ION:			
	- science - engineering	K. Anderson	Site property	warehouse	east of f	ormer retort	
ORILLED BY:			NORTHING:		EASTING:		
Zack Baile	y, Holocene Drilling		697981.38		1149583.06		
Sonic LAR			SURFACE ELEVATION: 2	252.1		COORDINATE SYSTEM: NAVD88	
DRILLING ME	THOD:		TOTAL DEPTH	(ft bgs):	DEPTH 1	TO WATER (ft bgs):	
5' x 4" Rod	ds		85		31		
SAMPLING M 4" Liner Ba	IETHOD/SAMPLER LENGTH: ag		BORING DIAME 4" inner/6" o		DRILL DATE: 10/2/2018		
Depth US0 (feet) Sym		scription and Observations R CONSTITUENT, odor, staining, sh	een, debris, etc.)	Drive/ Recovery	PID (ppm)	Sample ID	
84 — SN	Gray-brown, well-graded SA	G			0.0		

FLOVD	LCNIDED	PROJECT: West Coast Door	LOCATION: 31	02 S Cedar Tacoma WA	BORII	NG ID: SB-42
	ISNIDER	LOGGED BY:	BORING LOCAT			
strategy • sci	ience • engineering	K. Anderson		warehouse so	uth of	former retort
DRILLED BY:			NORTHING:	EASTING:		
Zack Bailey, Hold	ocene Drilling		697951.38		11495	53.06
DRILLING EQUIPMEN Sonic LAR	NT:		SURFACE ELEVATION: 2	E2 1	oordii NAVD	NATE SYSTEM: 88
DRILLING METHOD:			TOTAL DEPTH (•		O WATER (ft bgs):
5' x 4" Rods			80		32	
SAMPLING METHOD 4" Liner Bag	/SAMPLER LENGTH:		BORING DIAME 4" inner/6" ou	I	10/3/2	
Depth USCS (feet) Symbol	(color, texture, moisture, MAJOR	scription and Observations R CONSTITUENT, odor, stainin	g, sheen, debris, etc.)	Drive/ Recovery	PID (ppm)	Sample ID
	of CONCRETE. of rounded GRAVEL.					
Br At	rown, well-graded SAND with a second secon	brown. No odor.			0.0	
	17.5 feet bgs, becomes m					
ABBREVIATIONS:	20 100t bgg, becomes dry.		NOTES:			
	ground surface USCS = Unified String = denotes	Soil Classification System groundwater table	Recovered intervals evenly	y decompressed fo	or log un	less otherwise noted.

FI	OYDISNIDER	PROJECT: West Coast Door	LOCATION: 3	102 S Cedar , Tacoma WA	BORING	SB-42
	egy • science • engineering	LOGGED BY:	BORING LOCAT			
		K. Anderson		y warehouse so		rmer retort
DRILLEI Zack	D BY: Bailey, Holocene Drilling		NORTHING : 697951.38		ASTING: 1149553	3.06
DRILLIN Sonic	IG EQUIPMENT: LAR		SURFACE ELEVATION: 2	DE 2 1	oordina NAVD88	TE SYSTEM:
DRILLIN	IG METHOD:		TOTAL DEPTH			WATER (ft bgs):
5' x 4'	Rods		80		32	
	NG METHOD/SAMPLER LENGTH : er Bag		BORING DIAME 4" inner/6" o		RILL DAT 10/3/201	
Depth (feet)		scription and Observations R CONSTITUENT, odor, staining,	sheen, debris, etc.)		PID (ppm)	Sample ID
24 —	At 22 feet bgs, becomes mo Gray-brown, poorly-graded r SP Gray-brown, well-graded SA odor. Light rainbow sheet an	medium SAND with little ND with silt and gravel, i	moderate creosote		0.0	SB-42-25.5-26
28 —	Gray-brown, poorly-graded r	medium SAND , no odor o	or sheen.		0.1	SB-42-29-30
_					0.0	
32 —	Gray-brown, well-graded SA	ND with abundant grave	I and silt, no odor.		0.0	
36 — - 40	GRAVEL. Gray-brown, well-graded SA wet, no odor. Gray-brown, well-graded SA no odor.	ND with gravel and trace	e silt, moist to wet,		0.1	
	VIATIONS:		OTES:	ly docomproses of f	or loca umba -	o othornico neteri
rt bgs ppm =	= feet below ground surface USCS = Unified parts per million = denotes	Soil Classification System I	Recovered intervals even	iy decompressed fo	iog unies	s otherwise noted.

FLOYDISNIDER	PROJECT: West Coast Door	LOCATION: 310)2 S Cedar Tacoma WA	BORING	SB-42	
strategy • science • engineering	LOGGED BY:	BORING LOCATION	ON:			
strategy - scrence - engineering	K. Anderson	Site property	warehouse so	uth of fo	ormer retort	
DRILLED BY:		NORTHING:	EASTING:			
Zack Bailey, Holocene Drilling		697951.38		114955	3.06	
DRILLING EQUIPMENT: Sonic LAR		SURFACE ELEVATION: 25	TO 1	OORDINA NAVD8	ATE SYSTEM: 8	
DRILLING METHOD:		TOTAL DEPTH (ff	bgs): D	ЕРТН ТО	WATER (ft bgs):	
5' x 4" Rods		80		32		
SAMPLING METHOD/SAMPLER LENGTH:		BORING DIAMET	I	RILL DAT		
4" Liner Bag		4" inner/6" ou		10/3/20	18	
(feet) Symbol (color, texture, moisture, MAJOF	scription and Observations R CONSTITUENT, odor, staining, shee	n, debris, etc.)		PID ppm)	Sample ID	
Gray, poorly-graded medium SP At 44.5 feet bgs, sand becom SM Gray, poorly-graded coarse s Wood branch. SM Gray, poorly-graded coarse s Gray, poorly-graded coarse s Gray, firm sandy SILT, slight	nes coarse. silty SAND , wet, no odor.			0.0		
48 — ML Gray-brown, poorly-graded fi	ne SAND , moist, no odor.			0.0		
52 — Gray-brown, well-graded gra	velly SAND with silt, wet, n	o odor.		0.0		
At 60 feet bgs, silt content inc				0.0		
ABBREVIATIONS: ft bas = feet below ground surface LISCS = Unified 9	NOTE	ES: overed intervals eventy	decompressed for	r log unla	es athenvise noted	

FLOYDISNIDER	PROJECT: West Coast Door	LOCATION: 31	02 S Cedar , Tacoma WA	BORII	SB-42
strategy · science · engineering	LOGGED BY:	BORING LOCAT			
	K. Anderson		warehouse so		
DRILLED BY: Zack Bailey, Holocene Drilling		NORTHING : 697951.38	1	E ASTING 11495	
DRILLING EQUIPMENT:		SURFACE			NATE SYSTEM:
Sonic LAR		ELEVATION: 2	252.1	NAVD	
DRILLING METHOD:		TOTAL DEPTH (ft bgs):	DEPTH T	O WATER (ft bgs):
5' x 4" Rods		80		32	
SAMPLING METHOD/SAMPLER LENGTH: 4" Liner Bag		BORING DIAME 4" inner/6" or		10/3/2	
Depth USCS Soil Des (feet) Symbol (color, texture, moisture, MAJOI	scription and Observations R CONSTITUENT, odor, staining, sh	een, debris, etc.)	Drive/ Recovery	PID (ppm)	Sample ID
				0.0	
68 —				0.0	
At 70 feet bgs, silt content in	creases with depth.			0.0	
Gray, hard sandy SILT with r	ND with trace to little silt,			0.0	
At 75 feet bgs, sand coarsen 76 — SW. Bottom of boring = 80 feet bg					
ABBREVIATIONS:	NC	TES:			
ft bgs = feet below ground surface USCS = Unified :	Soil Classification System R groundwater table	ecovered intervals evenl	y decompressed f	or log un	less otherwise noted.

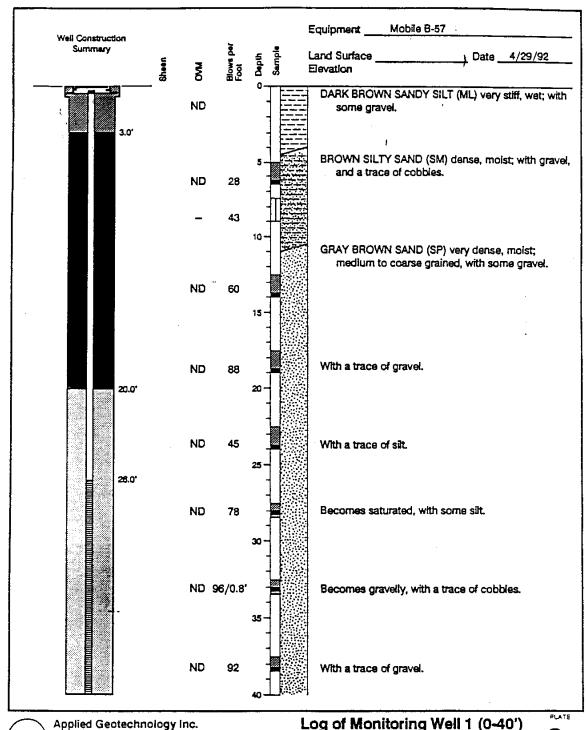
Strategy Science engineering Cocers Bry K. Anderson Side property warehouse west of former rotort	FT	OYDISNIDER	PROJECT: West Coast Door	LOCATION: 31	02 S Cedar Tacoma WA	BORING ID: SB-43
BRILLED BY: Zack Balley, Holocone Drilling BORILLING EXPORTED: SORICL LAR BORILLING METHOD: SAMPLER LENGTH: 4" Inner 16" Use 1" 104 Use			LOGGED BY:			
Zack Balley, Holocene Drilling BRILLING ECUIPMENT: SONIC LAR SURPACE			K. Anderson			
SORICLIAR BRILLING METHOD: 5 x 4" ROS SAMPLING METHOD: 5 x 4" ROS SORIC LOADER: 10 x 5 x 5 x 5 x 5 x 5 x 5 x 5 x 5 x 5 x						
DRILLING METHOD: 5 x 4" Rods 80 80 80 80 80 80 80 80 80 8				SURFACE ELEVATION: 2	E2 1	
SAMPLING METHODISAMPLER LENGTH: 4" Liner Bag 4" Inner 16" outer 4" Liner Bag 5 Clock Industrial Market Market Market Bag 6" of CONCRETE Dark brown, well-graded SAND with sitt and gravel, moist, no odor. At 2.5 feet bgs, drilled through concrete. At 8.5 feet bgs, pravel content increases. At 8.5 feet bgs, becomes brown and moist. Driver (Popular Sample ID) SW SM At 7 feet bgs, gravel content increases. At 8.5 feet bgs, becomes brown and moist. Driver (Popular Sample ID) Driver (Popular Sample ID) Driver (Popular Sample ID) Driver (Popular Sample ID) Sw SM At 7 feet bgs, gravel content increases. At 8.5 feet bgs, becomes brown and moist. Driver (Popular Sample ID) Driver (Popular	DRILLIN	IG METHOD:		TOTAL DEPTH (
A" LINET BIG. Depth USCS Soil Description and Observations Color. Instanting, sheen, stearts, etc.) Pipm Sample ID	5' x 4"	Rods		80		32
Color, touture, moletare, MALOR CONSTITUENT, odor, stanning, sheen, debris, etc.) Recovery (ppm) Sample ID					 	
Dark brown, well-graded SAND with silt and gravel, moist, no odor. Brick fragments from 0.5 to 2.5 feet bgs. At 2.5 feet bgs, drilled through concrete. At 2.5 feet bgs, drilled through concrete. At 3.5 feet bgs, drilled through concrete. At 8.5 feet bgs, becomes brown and moist. At 8.5 feet bgs, becomes brown and moist. Do 0 Brown, poorly-graded medium SAND with few fine rounded gavel, dry to slightly moist, no odor. SP At 15 feet bgs, very fine and dry powdery material, then brown, well-graded SAND with silt and gravel, moist. Gray-brown, poorly-graded medium SAND with gravel, moist. SP SM Gray-brown, poorly-graded medium SAND with gravel, moist. O 0 ABBREVIATIONS. Gray-brown, poorly-graded fine SAND with silt, moist, no odor. O 0 BO 15.5 B			scription and Observations R CONSTITUENT, odor, staining, sh	een, debris, etc.)		
Brick fragments from 0.5 to 2.5 feet bgs. At 2.5 feet bgs, drilled through concrete. At 2.5 feet bgs, drilled through concrete. At 7 feet bgs, gravel content increases. At 8.5 feet bgs, becomes brown and moist. O.0 SP Brown, poorly-graded medium SAND with few fine rounded gavel, dry to slightly moist, no odor. At 15 feet bgs, very fine and dry powdery material, then brown, well-graded SAND with silt and gravel, moist. Gray-brown, poorly-graded medium SAND with gravel, moist. Gray-brown, poorly-graded medium SAND with gravel, moist. O.0 ABBREVIATIONS: SP SM Gray-brown, poorly-graded fine SAND with silt, moist, no odor. O.0 NOTES: Recovered intervals evenly decompressed for log unless otherwise noted.	0					
SW-SM At 7 feet bgs, gravel content increases. At 8.5 feet bgs, becomes brown and moist. O.0 O.0 I2 — SP Brown, poorly-graded medium SAND with few fine rounded gavel, dry to slightly moist, no odor. At 15 feet bgs, very fine and dry powdery material, then brown, well-graded SAND with silt and gravel, moist. Gray-brown, poorly-graded medium SAND with gravel, moist. O.0 ABBREVIATIONS. It bgs - feet below ground surface USCS - Unified Soil Classification System NOTES: Recovered Intervals evenly decompressed for log unless otherwise noted.	4 —	Brick fragments from 0.5 to 2	2.5 feet bgs.	oist, no odor. -		0.0
Brown, poorly-graded medium SAND with few fine rounded gavel, dry to slightly moist, no odor. At 15 feet bgs, very fine and dry powdery material, then brown, well-graded SAND with silt and gravel, moist. Gray-brown, poorly-graded medium SAND with gravel, moist. O.0 Gray-brown, poorly-graded fine SAND with silt, moist, no odor. SP-SM ABBREVIATIONS: tt bgs = feet below ground surface USCS = Unified Soil Classification System NOTES: Recovered intervals evenly decompressed for log unless otherwise noted.		At 7 feet bgs, graver content		-		0.0
Brown, poorly-graded medium SAND with few fine rounded gavel, dry to slightly moist, no odor. At 15 feet bgs, very fine and dry powdery material, then brown, well-graded SAND with silt and gravel, moist. SW-SM Gray-brown, poorly-graded medium SAND with gravel, moist. O.0 Gray-brown, poorly-graded fine SAND with silt, moist, no odor. SP-SM Gray-brown, poorly-graded fine SAND with silt, moist, no odor. NOTES: Recovered intervals evenly decompressed for log unless otherwise noted.	_	At 8.5 feet bgs, becomes bro	own and moist.			
well-graded SAND with silt and gravel, moist. Gray-brown, poorly-graded medium SAND with gravel, moist. O.0 Gray-brown, poorly-graded fine SAND with silt, moist, no odor. SP-SM 20 ABBREVIATIONS: It bgs = feet below ground surface USCS = Unified Soil Classification System NOTES: Recovered intervals evenly decompressed for log unless otherwise noted.	- 12 —	Michigan I.	m SAND with few fine rou	ınded gavel, dry	-	0.0
Gray-brown, poorly-graded fine SAND with silt, moist, no odor. SP-SM 20 ABBREVIATIONS: ft bgs = feet below ground surface USCS = Unified Soil Classification System NOTES: Recovered intervals evenly decompressed for log unless otherwise noted.	16 —	well-graded SAND with silt a	nd gravel, moist.			0.0
ABBREVIATIONS: ft bgs = feet below ground surface USCS = Unified Soil Classification System NOTES: Recovered intervals evenly decompressed for log unless otherwise noted.	_	Gray-brown, poorly-graded f	ne SAND with silt, moist,	no odor.		0.0
ft bgs = feet below ground surface USCS = Unified Soil Classification System Recovered intervals evenly decompressed for log unless otherwise noted.		INC. COLO	NO	TEC:		
	ft bgs	= feet below ground surface USCS = Unified	Soil Classification System Re		decompressed for	or log unless otherwise noted.

EL OVD I CNIDED	PROJECT: West Coast Door	LOCATION: 310	2 S Cedar Tacoma WA	BORING ID: SB-43		
FLOYDISNIDER		ORING LOCATION:				
strategy • science • engineering	varehouse west of former retort					
DRILLED BY:		NORTHING:				
Zack Bailey, Holocene Drilling		697981.38		1149523.06		
DRILLING EQUIPMENT: Sonic LAR		SURFACE ELEVATION: 25	: 1 · · · · · · · · · · · · · · · · · ·	OORDINATE SYSTEM: NAVD88		
DRILLING METHOD:		TOTAL DEPTH (ft	bgs): D	EPTH TO WATER (ft bgs):		
5' x 4" Rods		80		32		
SAMPLING METHOD/SAMPLER LENGTH: 4" Liner Bag		4" inner/6" out		PRILL DATE: 10/4/2018		
(feet) Symbol (color, texture, moisture, MAJOR	scription and Observations R CONSTITUENT, odor, staining, shee	n, debris, etc.)	Drive/ Recovery (PID (ppm) Sample ID		
20 Brown, well-graded SAND w	ith silt and gravel, moist.					
Gray-brown, poorly-graded fi	ne SAND with silt, moist, n	o odor.		0.0		
At 22.5 feet bgs, becomes gr	ay with a slight odor. No sh	neen.		0.0		
Gray, well-graded SAND with Slight rainbow sheet and clear				2.5		
Gray, poorly-graded fine SAI and film and trace oily produce	Gray, poorly-graded fine SAND , creosote odor. Trace rainbow sheen and film and trace oily product.					
At 30 feet bgs, becomes ligh				394.0 SB-43-31-31.5		
and trace brown oily film.	and trace brown oily film.					
36 — odor.						
Gray, poorly-graded medium rainbow sheen and brown oil		dor. Light		SB-43-37.5-38		
SP.				13.3		
40 [WASSER] ABBREVIATIONS:	NOT	L ES:				
ft bgs = feet below ground surface USCS = Unified S			decompressed fo	or log unless otherwise noted.		
ppm = pans per million	groundwater table					

FLOYDISNIDER	PROJECT: West Coast Door	LOCATION: 31	02 S Cedar , Tacoma WA	BORIN	SB-43	
		BORING LOCAT				
strategy • science • engineering	K. Anderson		warehouse we	est of fo	ormer retort	
DRILLED BY:		NORTHING:		EASTING:		
Zack Bailey, Holocene Drilling		697981.38		114952	23.06	
DRILLING EQUIPMENT:		SURFACE) [2 1		IATE SYSTEM:	
Sonic LAR		ELEVATION: 2	252.1	NAVD8	38	
DRILLING METHOD:		TOTAL DEPTH (O WATER (ft bgs):	
5' x 4" Rods		80		32		
SAMPLING METHOD/SAMPLER LENGTH: 4" Liner Bag		BORING DIAME 4" inner/6" o	 	10/4/20		
		4 111110170 0	Drive/		0.10	
	escription and Observations OR CONSTITUENT, odor, staining, she	en, debris, etc.)		PID (ppm)	Sample ID	
40						
Gray, well-graded SAND wi	th silt and gravel, wet, no oc	dor.		0.1		
				-	SB-43-41-41.5	
At 42 fact has silt content in	ncreases and sand coarsen	c with donth	-			
At 42 feet bys, sin content i	nicreases and sand coarsen	s willi depili.				
l				0.1		
44 —						
				0.1		
Gray, firm SILT , no odor.			-	-	SB-43-47-47.5	
				0.0		
48						
Gray, well-graded very fine	SAND, moist to wet.					
Gray, well-graded SAND w	ith silt and gravel, wet, no o	dor.				
At 51 feet bgs, becomes we	et.					
52 —						
				0.0		
At 55 feet bgs, becomes gr	ay-brown.					
56						
30 7						
				0.0		
At 57 feet bgs, becomes br	own.					
(1) (1) (1) (1) (1) (1) (1) (1) (1) (1)						
[
60 At 60 feet bgs, becomes de	ense and silt content increas	es				
ABBREVIATIONS:	NOT	ES:				
ft has - feet helow around surface IISCS - Unified	NOT Classification System RAC	covered intervals event	v decompressed to	or inditini	ess otherwise noted	

FLOYDISNIDER	PROJECT: West Coast Door	LOCATION: 310	02 S Cedar Tacoma WA	BORIN	SB-43
strategy • science • engineering	LOGGED BY:	BORING LOCATI			
strategy • science • engineering	K. Anderson		warehouse we	est of fo	rmer retort
DRILLED BY:		NORTHING:		ASTING	
Zack Bailey, Holocene Drilling		697981.38		114952	23.06
DRILLING EQUIPMENT: Sonic LAR		SURFACE ELEVATION: 2	E 2 1	<mark>oordin</mark> NAVD8	ATE SYSTEM: 88
DRILLING METHOD:		TOTAL DEPTH (f	t bgs): D	EPTH TO	O WATER (ft bgs):
5' x 4" Rods		80		32	
SAMPLING METHOD/SAMPLER LENGTH:		BORING DIAMET	l l	RILL DA	
4" Liner Bag		4" inner/6" ou		10/4/20)18
Depth USCS Soil Des (feet) Symbol (color, texture, moisture, MAJOR	scription and Observations R CONSTITUENT, odor, staining, sheen,	debris, etc.)	Drive/ Recovery (PID (ppm)	Sample ID
64 —				0.0	
At 65.5 feet bgs, silt content		-		0.0	
Gray-brown, very fine poorly	-graded SAND , moist to wet,	no odor.			SB-43-71-71.5
Gray-brown, dense silty SAN				0.0	
76 — Gray, hard sandy SILT with o	gravel, dry to slightly moist, n	o odor.		0.0	
80 SM Gray, silty SAND . Bottom of	boring = 80 feet bgs.				
ABBREVIATIONS:	NOTES Recovered		decompressed for	or log upla	ess otherwise noted

ft bgs = feet below ground surface USCS = Unified Soil Classification System ppm = parts per million = denotes groundwater table



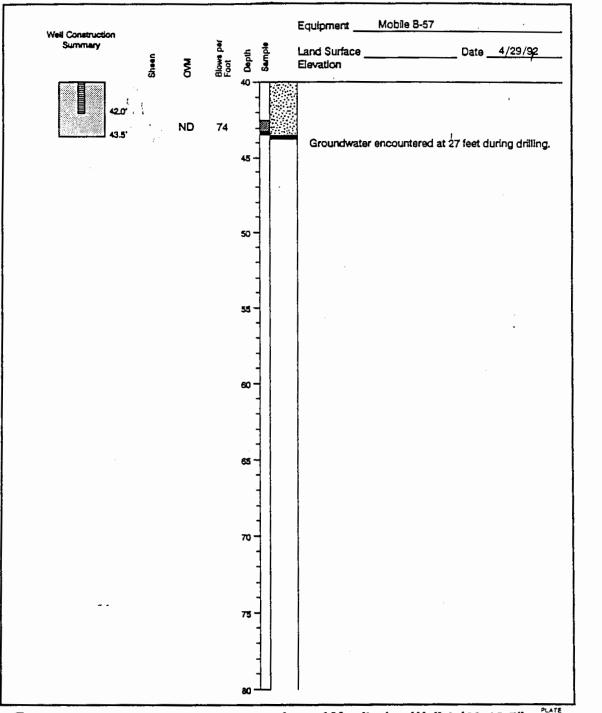
Geotechnical Engineering Geology & Hydrogeology

Log of Monitoring Well 1 (0-40')

Puget Sound National Bank/West Coast Door - Phase II 3a Tacoma Washington

DATE REVISED / JOB NUMBER 15,536.003 SEŞ JUL 18 Jun 92

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Applied Geotechnology Inc. Geotechnical Engineering Geology & Hydrogeology

Log of Monitoring Well 1 (40-43.5')
Puget Sound National Bank/West Coast Door - Phase II 3b
Tacoma, Washington

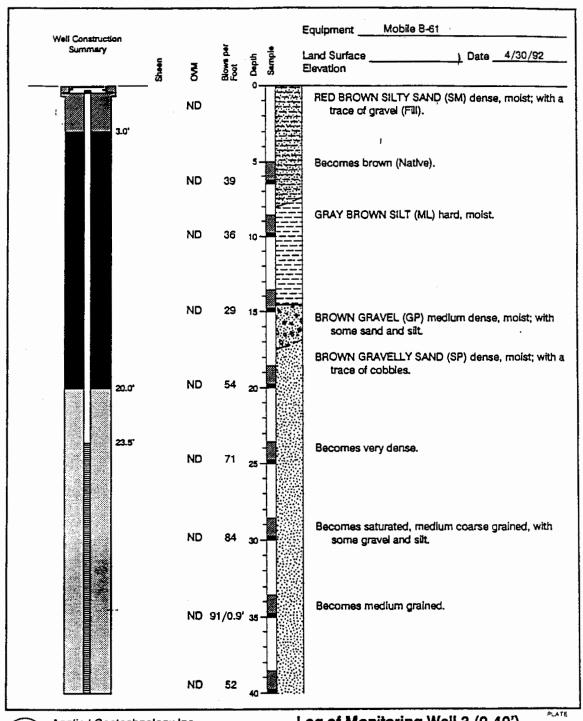
JOS NUMBER 15,536,003

ORAWN SES

DATE 18 Jun 92

REVISED

DATE /

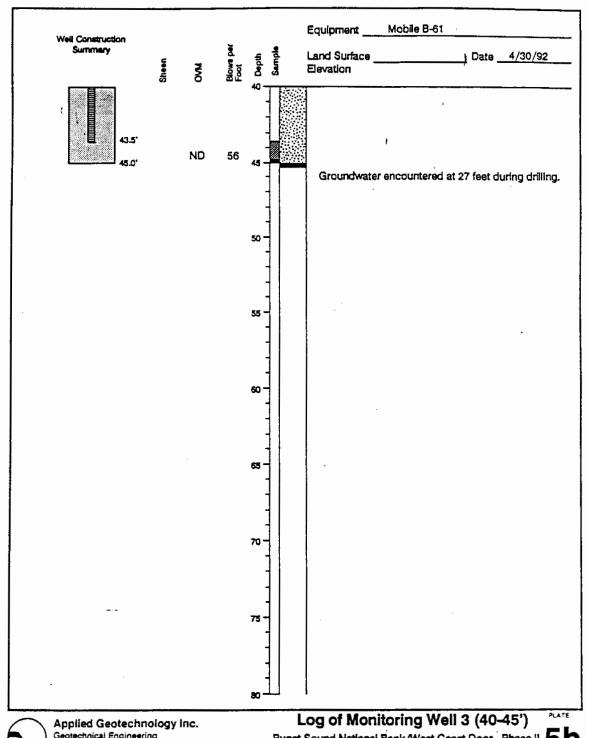


Applied Geotechnology Inc. Geotechnical Engineering Geology & Hydrogeology

Log of Monitoring Well 3 (0-40')

Puget Sound National Bank/West Coast Door - Phase II 5a Tacoma, Washington

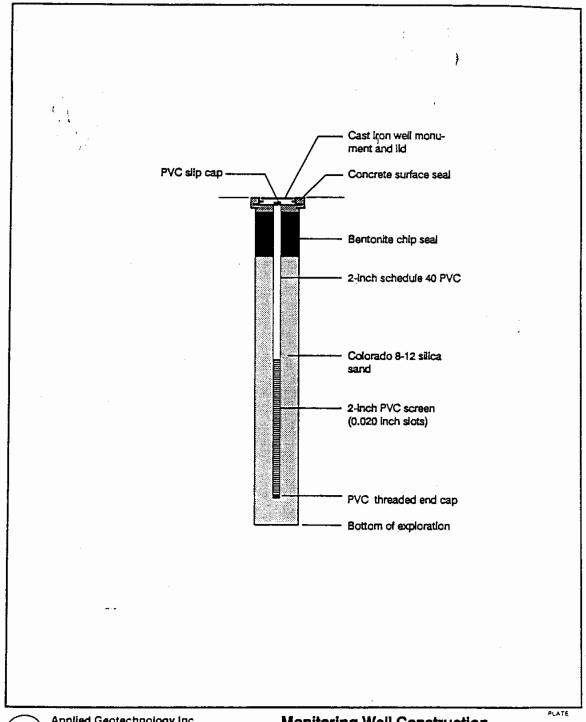
DATE PEVISED / CATE 18 Jun 92 15,536.003 SES



Geotechnical Engineering Geology & Hydrogeology

Puget Sound National Bank/West Coast Door - Phase || 5b Tacoma, Washington

JOB NUMBER 15,536,003 DRAWN APPROVED THE DATE 18 Jun 92 REVISED ! DATE





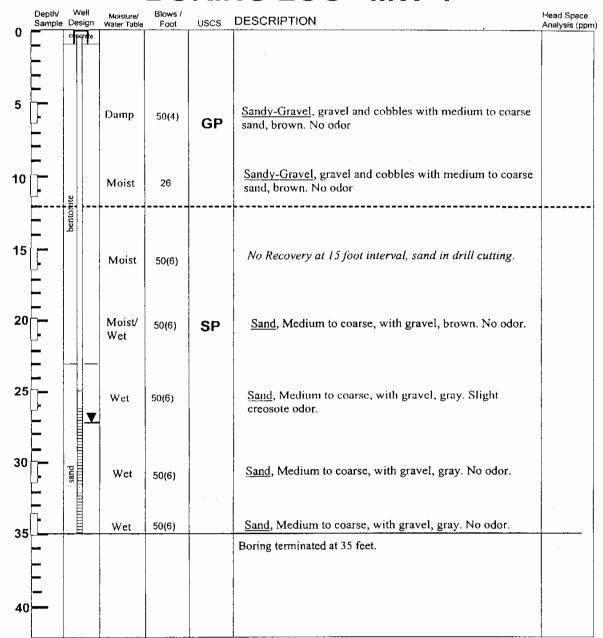
Applied Geotechnology Inc. Geotechnical Engineering Geology & Hydrogeology

Monitoring Well Construction
Puget Sound National Bank/West Coast Door - Phase II
Tacoma, Washington

2

JOB NUMBER DRAWN APPROVED CATE REVISED DATE
15.536.003 SES 18 Jun 92

BORING LOG MW-4



Sampler: Split-spoon. Soil sampled collected every 5 feet beginning at 2.5 feet below ground surface. Driller: Cascade Drilling, Inc. Hollow-stem auger.

Monitoring Well: 2"-PVC, Screened 25' to 35', 0.010" Slot



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ENVIRONMENTAL ASSOCIATES, INC.

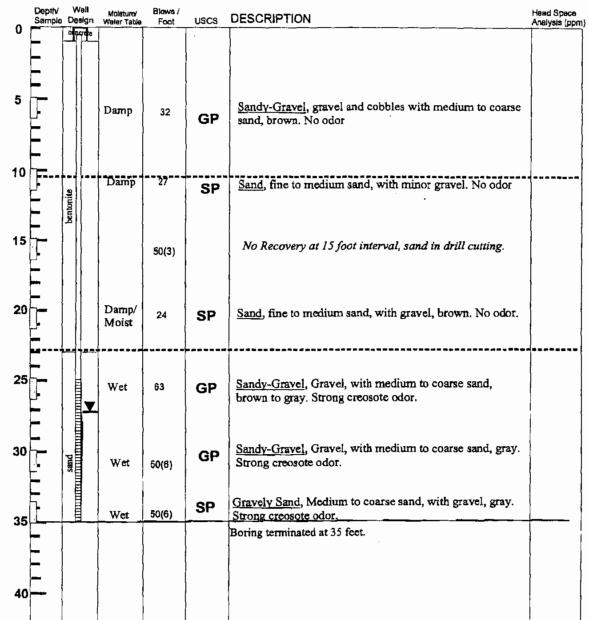
1380 - 112th Avenue NE, Suite 300 Bellevue, Washington 98004

BORING LOG MW-4

West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Date:	Logged by:	Plato:
07/28/06	RBR	A-5
		** '

BORING LOG MW-5



Sampler: Split-spoon. Soll sampled collected every 5 feet beginning at 2.5 feet below ground surface. Driller: Cascade Drilling, Inc. Hollow-stern auger.

Monitoring Well: 2"-PVC, Screened 25' to 35', 0.010" Slot



ENVIRONMENTAL ASSOCIATES, INC.

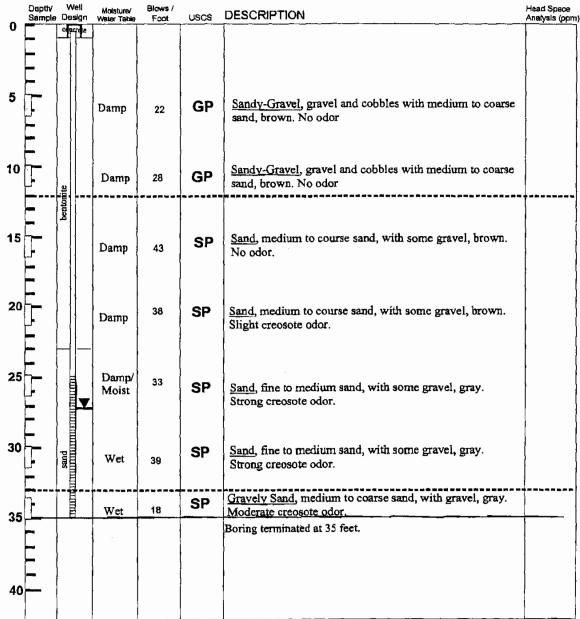
1380 - 112th Avenue NE, Suite 300 Beilevue, Washington 98004

BORING LOG MW-5

Former West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number:	Date:	Logged by:	Plate:
JN-26192-3	09/18/06	RBR	A-6

BORING LOG MW-6



Sampler: Spilt-spoon. Soil sampled collected every 5 feet beginning at 2.5 feet below ground surface. Driller: Cascade Drilling, Inc. Hollow-stem auger.

Monitoring Well: 2"-PVC, Screened 25' to 35', 0.010" Slot



1380 - 112th Avenue NE, Suite 300 Beilevne, Washington 98004

BORING LOG MW-6

Former West Coast Door Property 3133 South Cedar Street Tacoma, Washington 98409

Job Number:	Date:	Logged by:	Plate:
JN-26192-3	09/18/06	RBR	A-7

Date/Time Started : 1-26-07/0827 Date/Time Completed: 1-26-07/1415 LOG OF WELL MW-7 **Total Boring Depth** : 52.5' Total Well Depth . 40 (Page 1 of 3) Depth to water ATD : 27.5 Elevation (ft) : NA Site Name: West Coast Door Property **Drilling Method** : HSA : D+M S.S. 3" diameter 18' Client: William Swensen Sampler Type Drive Hammer (lbs) : 140 PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 Project #: 112-001 Depth In Feet Well: MW-7 Recovery Blow Count PID (ppm) Graphic Description Sample ID 0.0 - 0.5 Concrete. Concrete 0.5 - 1.0 Pea gravel. GP Surface Seal 1.0 - 2.5 Silty SAND with strong creosote odor, SM 2.5 - 3.0 Concrete. 5. 5.0 - 6.5 Cobble. 50/3" 7.5 - 9.0 Broken Cobble. 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\WW7.bo 50/1" Beptonitelank Casing Annular Seal 10-10.0 - 11.5 SAND with gravel, trace silt (80% medium to coarse sand, 15% fine to coarse SP 50/2" gravel, 5% silt), brown, moist. 12.5 - 14.0 GRAVEL, trace sand, trace silt (90% fine gravel, 5% coarse sand, 5% silt), brown, GP 50/3" 44.8 moist. 15-15.0 - 16.5 GRAVEL, trace sand, trace silt (90% fine to coarse gravel, 5% coarse sand, 5% silt), GP 50/3" 30 brown, moist. 17.5 - 18.25 Gravelly SAND trace silt (55% SP coarse sand, 40% fine gravel, 5% silt), brown, 67.0 MW7-17.5-19 50 50/6" moist, creosote odor. SP 18.25 - 19.0 SAND (100% fine sand), brown, moist, strong creosote odor. Drilling Company : Cascade Drilling, Inc. Drilling Foreman : Steve Choate LOG OF WELL MW-7 Equipment : CME 65 Pacific Crest Rep. : Annica Nord (Page 1 of 3)

Date/Time Started : 1-26-07/0827 Date/Time Completed: 1-26-07/1415 LOG OF WELL MW-7 Total Boring Depth : 52.5' Total Well Depth . 40 (Page 2 of 3) Depth to water ATD : 27.5 Elevation (ft) : NA Site Name: West Coast Door Property **Drilling Method** : HSA Client: William Swensen Sampler Type : D+M S.S. 3" diameter 18' Drive Hammer (lbs) : 140 PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 N D R T H B E N D , W A 9 B D 4 5 Project #: 112-001 Depth In Feet Recovery Well: MW-7 Blow Count (mdd) Samples Graphic Description USCS PID (Sample ID 20 20.0 - 21.5 Gravelly SAND trace silt (55% coarse sand, 40% fine gravel, 5% silt), brown, moist, SP 50/3" 64 Bentonite creosote odor. Annular Seal 2" PVC Blank Casing 22.5 - 24.0 SAND trace gravel (95% medium to coarse sand, 5% coarse gravel), gray, moist, SP 20 50/4" 60.2 creosote odor. 25 25.0 - 26.5 SAND trace gravel (95% medium to coarse sand, 5% coarse gravel), gray, moist, SP 50/6" MW7-25-26.5 65 creosote odor. ▼ 27.5 - 29.0 SAND (100% fine to coarse sand), 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\WW7.bo gray, wet, slight creosote odor. SP 100 29/30/39 22 30. 30.0 - 31.5 SAND (100% fine to coarse sand), gray, wet, slight creosote odor. SP 33/36/39 32.8 -#2/12 Sand Pack 2" PVC 32.5 - 34.0 GRAVEL minor sand (85% fine to 0.010 slot screen coarse gravel, 15% medium to coarse sand), GP 27/30/32 61.2 dark gray, wet, slight musty odor. 35 35.0 - 36.5 NO RECOVERY 29/31/32 37.5 - 39.0 NO RECOVERY. 50/2" **Drilling Company** : Cascade Drilling, Inc. Drilling Foreman : Steve Choate LOG OF WELL MW-7 Equipment : CME 65 Pacific Crest Rep. : Annica Nord (Page 2 of 3)

Date/Time Started : 1-26-07/0827 Date/Time Completed: 1-26-07/1415 LOG OF WELL MW-7 Total Boring Depth : 52.5' Total Well Depth : 40' (Page 3 of 3) Depth to water ATD : 27.5 Elevation (ft) : NA Site Name: West Coast Door Property **Drilling Method** : HSA Client: William Swensen Sampler Type : D+M S.S. 3" diameter 18' Drive Hammer (lbs) : 140 PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 N D R T H B E N D , W A 9 B D 4 5 Project #: 112-001 Depth In Feet Recovery Well: MW-7 Blow Count (mdd) Samples Graphic Description USCS PID (Sample ID % 40 40.0 - 41.0 SAND (100% fine to medium sand), SP gray, wet, slight creosote odor. 20 50/4" 22 41.0 - 41.5 Sandy GRAVEL (50% fine to coarse GP gravel, 50% medium to coarse sand), gray, wet, slight creosote odor. 42.5 - 44.0 Gravelly SAND (60% fine to coarse sand, 40% fine to coarse gravel), gray, wet, SP 35 50/6" 44.8 creosote odor. 45 45.0 - 46.5 SAND (100% fine to coarse sand), gray, wet, creosote odor. SP 32-50/4" 34.8 MW7-GW-45 Bentonite 47.5 - 49.0 NO RECOVERY. 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\WW7.bo 50/6" 50 50.0 - 51.5 SAND (100% fine to coarse sand), gray, wet, heavy rainbow sheen, strong SP 37.1 100 50/4" petroleum odor. 52.5 - 54.0 NO RECOVERY. MW7-52.5-GW 55 60 **Drilling Company** : Cascade Drilling, Inc. Drilling Foreman : Steve Choate LOG OF WELL MW-7 Equipment : CME 65 Pacific Crest Rep. : Annica Nord (Page 3 of 3)

Date/Time Started : 1-31-07/0850 Date/Time Completed: 1-31-07/1100 LOG OF WELL MW-8 **Total Boring Depth** : 40' Total Well Depth : 40' (Page 1 of 2) Depth to water ATD : 30' Elevation (ft) : NA Site Name: West Coast Door Property **Drilling Method** : HSA Sampler Type : D+M S.S. 2" diameter 18' Client: William Swensen Drive Hammer (lbs) : 140 Project #: 112-001 Depth In Feet Recovery Well: MW-8 Blow Count PID (ppm) Graphic Description Sample ID 0.0 - 0.5 Concrete. Concrete 0.5 - 5.0 Soil cuttings are sandy with mainly Surface Seal rounded cobbles. GP 5 5.0 - 6.5 NO RECOVERY. 50/1" Driller comments that below 8 feet, he is not drilling in fill material. 08-13-2008 NPacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\MW8.bor 10.0 - 11.5 Sandy GRAVEL with silt, fill, dry. GΡ 50/2" 32.7 2" PVC Blank Casing Bentonite Annular Seal 15 15.0 - 16.5 SAND minor silt, minor gravel (75% fine to coarse sand, 15% silt, 10% fine gravel), 166 SP 50/4" MW8-15-16.5 30 brown, dry. 20 20.0 - 21.5 SAND minor gravel, trace silt (85% fine to coarse sand, 10% fine gravel, 5% silt), SP 20 50/5" 96.6 brown-yellow, slightly moist. **Drilling Company** : Cascade Drilling, Inc. Drilling Foreman : Steve Choate LOG OF WELL MW-8 Equipment : CME 65 Pacific Crest Rep. : Annica Nord (Page 1 of 2)

Date/Time Started : 1-31-07/0850 Date/Time Completed: 1-31-07/1100 LOG OF WELL MW-8 **Total Boring Depth** : 40' Total Well Depth : 40' (Page 2 of 2) Depth to water ATD : 30' Elevation (ft) : NA Site Name: West Coast Door Property **Drilling Method** : HSA Sampler Type : D+M S.S. 2" diameter 18' Client: William Swensen Drive Hammer (lbs) : 140 PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 N D R T H B E N D , W A 9 B D 4 5 Project #: 112-001 Depth In Feet Well: MW-8 Recovery **Blow Count** PID (ppm) Samples Graphic uscs Description Sample ID 22 Bentonite Annular Seal 2" PVC Blank Casing 25.0 - 26.5 SAND trace gravel, trace silt (90% fine to coarse sand, 5% fine gravel, 5% silt), SP 50/6" 69/115 MW8-25-26.5 50 brown, slightly moist. 27 ▼ 30.0 - 31.5 SAND trace silt (95% fine to coarse 08-13-2008 NPacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\MW8.bor sand, 5% silt), brown, wet, no odor. SP 50 50/6" 35.9 -#2/12 Sand Pack 32 2" PVC 0.010 slot screen 35.0 - 36.5 SAND minor gravel, trace silt (85% fine to coarse sand, 10% fine gravel, 5% silt), SP 70 50/6" 93.6 brown-gray, wet, assorted cobbles. 37 40.0 - 41.5 SAND (100% fine to coarse sand), gray-brown, wet, no odor. SP 55 50/6" 23 42 **Drilling Company** : Cascade Drilling, Inc. Drilling Foreman : Steve Choate LOG OF WELL MW-8 Equipment : CME 65 Pacific Crest Rep. : Annica Nord (Page 2 of 2)

		LOG OF W		age 1 of	4)			Date/Time Sta Date/Time Co Total Boring D Total Well De Depth to wate	mpleted epth oth	: 9-6-2007/0915 : 9-7-2007/1100 : 70' : 70' : 25'
			Site Name: West C	oast Doc	r Pro	opert	у	Elevation Drilling Metho		: 245.99' : Sonic
15		ROLL TO SE	Client: Willia	am Swens	sen			Sampler Type		: Sonic Core Sampler
EEE!	BEN	DIGG BOULEVARD NORTH PO BOX 952 H B E N D , W A 9 B D 4 5	Project #:	: 112-001						
Depth In Feet	Samples	Description		nscs	Graphic	% Recovery	PID (ppm)	Lab No.	Well:	MW-9
0-		0.0 - 0.5 Asphalt. 0.5 - 2.5 GRAVEL, minor sand, trace s coarse gravel, 10% fine to coarse sand brown-gray, moist, no odor.	llt, (85% fine to l, 5% silt),	GP		100	0.0			—Concrete Surface Seal
-		2.5 - 5.0 GRAVEL, minor sand, trace s coarse gravel, 10% fine to coarse sand moist, no odor.	ilt, (85% fine to l, 5% silt), brown,	GP		100	0.0	-		
5— - - -	\bigvee	5.0 - 7.5 GRAVEL, minor sand, trace s coarse gravel, 10% fine to coarse sand moist, no odor.		GP		100	-	-		
-		7.5 - 10.0 GRAVEL, minor sand, trace coarse gravel, 10% fine to coarse sand moist, no odor.	silt, (85% fine to I, 5% silt), brown,	GP		100	0.0	-		4" PVC Blank Casi
10	\bigvee	10.0 - 12.0 SAND, (100% fine to medit moist, no odor. 12.0 - 12.5 Sandy GRAVEL, (65% fine		SP		100	0.0	-		Bentonite Annular Seal
		12.5 - 12.5 Sandy GRAVEL, (65 % line to coarse sand), light brown-c 12.5 - 15.0 Sandy GRAVEL, (65 % fine 35 % fine to coarse sand), light brown-c	ray, moist, no odorto coarse gravel,	GP		100	-	-		
15 — - - -	X	15.0 - 17.5 Gravelly SAND, trace silt, (sand, 40% fine to coarse gravel, 5% si moist, no odor.	55% fine to coarse t), brown-gray,	SP		100	-	-		
- - - 20-		17.5 - 20.0 Gravelly SAND, trace silt, (sand, 40% fine to coarse gravel, 5% si moist, no odor.	55% fine to coarse t), brown-gray,	SP		100	1.5	-		
Drilli		ompany : Boat Longyear Drilling, Inc.								
Equi	pme							LOG	G OF \	WELL MW-9
Pacif	fic C	rest Rep. : Annica Nord								(Page 1 of 4)

Date/Time Started : 9-6-2007/0915 Date/Time Completed: 9-7-2007/1100 LOG OF WELL MW-9 **Total Boring Depth** : 70' Total Well Depth . 70' (Page 2 of 4) Depth to water ATD : 25' Elevation : 245.99' Site Name: West Coast Door Property **Drilling Method** : Sonic Sampler Type : Sonic Core Sampler Client: William Swensen PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 N D R T H B E N D , W A 9 B D 4 5 Project #: 112-001 In Feet Recovery Well: MW-9 (mdd) Samples Graphic Depth I Description PID (Lab No. % 20 20.0 - 21.0 SAND, minor gravel, trace silt (85% fine to SP 100 0.0 coarse sand, 10% fine gravel, 5% silt), dark brown, moist, 21.0 - 23.0 SAND, with silt (80% fine to coarse sand, 20% silt), gray-brown, moist, no odor. SP 100 0.0 23.0 - 24.0 GRAVEL, with sand, trace silt, (75% fine to GP coarse gravel, 20% fine to coarse sand, 5% silt), gray, 100 MW9-24-25 3.8 moist, no odor. SP-SM 24.0 - 25.0 Silty SAND, (60% fine sand, 40% silt), gray, wet, ▼ 25 creosote odor, assorted cobbles. 25.0 - 27.0 SAND, (100% find sand), gray, wet, strong creosote odor. SP 100 0.0 27.0 - 27.5 GRAVEL, with sand, (80% fine to coarse gravel, GP 20% fine to coarse sand), gray, wet, rainbow sheen, strong 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\WW9.bo creosote odor. -4" PVC Blank Casing 27.5 - 30.0 GRAVEL, with sand, (80% fine to coarse gravel, GP 100 0.0 20% fine to coarse sand), gray, wet, heavy rainbow sheen, strong creosote odor. 30 Bentonite Annular 30.0 - 32.5 GRAVEL, with sand, (80% fine to coarse gravel, Seal 20% fine to coarse sand), gray, wet, heavy rainbow sheen, strong creosote odor. GP 100 2.6 32.5 - 35.0 Gravelly SAND, (55% fine to coarse sand, 45% fine to coarse gravel), gray, wet, strong creosote odor, heavy rainbow sheen. SP 100 1.6 35 35.0 - 37.0 SAND, minor gravel, (90% fine to coarse sand, 10% fine to coarse gravel), gray, wet, creosote odor. SP 100 0.0 37.0 - 37.5 GRAVEL, with sand, (75% fine to coarse gravel, GP 25% fine to coarse sand), gray, wet, creosote odor, slight sheen evident on soil. 37.5 - 40.0 GRAVEL, with sand, (75% fine to coarse gravel, GΡ 100 MW9-40-GW 25% fine to coarse sand), gray, wet, creosote odor, slight sheen evident on soil. 40 **Drilling Company** : Boat Longyear Drilling, Inc. **Drilling Foreman** : Dale LOG OF WELL MW-9 Equipment : Sonic Pacific Crest Rep. : Annica Nord (Page 2 of 4)

Date/Time Started : 9-6-2007/0915 Date/Time Completed: 9-7-2007/1100 LOG OF WELL MW-9 **Total Boring Depth** : 70' Total Well Depth . 70' (Page 3 of 4) Depth to water ATD : 25' Elevation : 245.99' Site Name: West Coast Door Property **Drilling Method** : Sonic Sampler Type : Sonic Core Sampler Client: William Swensen PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 N D R T H B E N D , W A 9 B D 4 5 Project #: 112-001 In Feet Recovery Well: MW-9 (mdd) Samples Graphic Depth I Description PID (Lab No. % 40 40.0 - 42.5 SAND (100% fine to medium sand), gray, wet, creosote odor. SP 100 42.5 - 44.5 SAND (100% fine to medium sand), gray, wet, creosote odor. SP 100 44.5 - 47.0 Sandy SILT, (50% silt, 50% fine sand), gray, 45 moist, very dense, no odor. ML100 0.0 MW9-45-47 47.0 - 50.0 Sandy GRAVEL, trace silt, (55% fine to coarse gravel, 40% fine to coarse sand, 5% silt), brown, wet, no NPacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\MW9.bo -4" PVC Blank Casing GP 100 0.0 50.0 - 52.5 GRAVEL, with sand, trace silt, (75% fine to coarse gravel, 20% fine to coarse sand, 5% silt), brown, Bentonite Annular wet, no odor. Seal 52.5 - 53.5 GRAVEL, with sand, trace silt, (75% fine to 50 coarse gravel, 20% fine to coarse sand, 5% silt), brown, wet, no odor. 53.5 - 54.0 GRAVEL, trace sand, trace silt, (90% fine to GΡ 100 coarse gravel, 5% fine to coarse sand, 5% silt), brown, wet, no odor. 54.0 - 55.0 Sandy GRAVEL, trace silt, (55% fine to coarse GP gravel, 40% fine to coarse sand, 5% silt), brown, wet, creosote odor. GP 100 55.0 - 55.5 GRAVEL, trace sand, trace silt, (90% fine to coarse gravel, 5% fine to coarse sand, 5% silt), brown, GP wet, slight creosote odor. 55 55.5 - 57.5 Gravelly SAND, (55% fine to coarse sand, 45% GP fine to coarse gravel), gray-brown, wet, very sight to no creosote odor. 100 0.0 MW9-55.5-57.5 SP 57.5 - 59.5 Sandy GRAVEL, trace silt, (70% fine to coarse gravel, 25% fine to coarse sand, 5% silt), wet, brown, slight 59.5 - 60.0 GRAVEL, trace sand, trace silt, (90% fine to coarse gravel, 5% fine to coarse sand, 5% silt), brown, GΡ 100 wet, no odor. #2/12 Sand Pack GP 60 : Boat Longyear Drilling, Inc. **Drilling Company** Drilling Foreman : Dale LOG OF WELL MW-9 Equipment : Sonic Pacific Crest Rep. : Annica Nord (Page 3 of 4)

Date/Time Started : 9-6-2007/0915 Date/Time Completed: 9-7-2007/1100 LOG OF WELL MW-9 **Total Boring Depth** : 70' Total Well Depth : 70' (Page 4 of 4) Depth to water ATD : 25' Elevation : 245.99' Site Name: West Coast Door Property **Drilling Method** : Sonic Sampler Type Client: William Swensen : Sonic Core Sampler PACIFIC CREST ENVIRONMENTAL 1533 BENDIGO BOULEVARD NORTH PO BOX 952 N D R T H B E N D , W A 9 B D 4 5 Project #: 112-001 Depth In Feet Well: MW-9 Recovery PID (ppm) Samples Graphic Description Lab No. 60 60.0 - 64.0 GRAVEL, minor sand, trace silt, (85% fine to coarse gravel, 10% fine to coarse sand, 5% silt), brown, wet, no odor. GΡ 100 4" PVC 0.010 slot screen 64.0 - 67.0 GRAVEL, minor sand, trace silt, (85% fine to coarse gravel, 10% fine to coarse sand, 5% silt), brown, 65 -#2/12 Sand Pack wet, no odor. GP 100 0.0 67.0 - 67.5 GRAVEL, minor sand, trace silt, (85% fine to GP coarse gravel, 10% fine to coarse sand, 5% silt), brown, 08-13-2008 \\Pacific-8e185af\public\Project Files\112 Swensen\112-001 West Coast Door Property\Boring Logs\WW9.bo wet, no odor. 100 0.0 MW9-68-70 67.5 - 70.0 Sandy SILT, (60% silt, 40% very fine sand), ML gray, moist, no odor, very dense. 70 75 80-**Drilling Company** : Boat Longyear Drilling, Inc. Drilling Foreman : Dale LOG OF WELL MW-9 Equipment : Sonic Pacific Crest Rep. : Annica Nord (Page 4 of 4)

Ground Surf Elev. & Datum: 244.5 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 697945.61 Longitude/Easting: 1149387.71 Casing Elevation: 244.22 ft Monitoring Well ID: MW-10

Drill Date: June 10, 2010 **Logged By:** John LaManna

Drilled By: Curtis Askew / Cascade Drilling

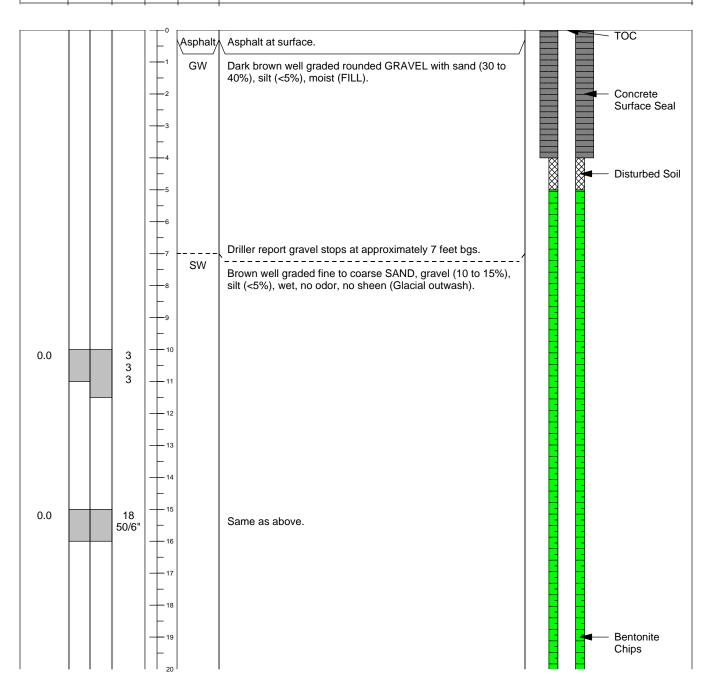
Drill Type: CME 75; 4-inch HSA
Sample Method: 18" D&M Split-spoon
Boring Diameter: 8"
Client: Bill Swensen
Project: Swensen-WCD
Site Location:3133 Cedar St,

Boring Depth (ft bgs): 46 ft bgs **Groundwater ATD (ft bgs):** 21.84* Tacoma, WA.

Remarks: *Depth to water from top of casing from 6/22/10 groundwater sampling.

PID Reading DRIVE / BLOW DEPTH USCS SOIL DESCRIPTION AND OBSERVATIONS: (color, texture, sample ID RECOVERY COUNT FT BGS SYMBOL moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)

MONITORING WELL DETAIL



Ground Surf Elev. & Datum: 244.5 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 697945.61 Longitude/Easting: 1149387.71 Casing Elevation: 244.22 ft Monitoring Well ID: MW-10

Drill Date: June 10, 2010 **Logged By:** John LaManna

Drilled By: Curtis Askew / Cascade Drilling

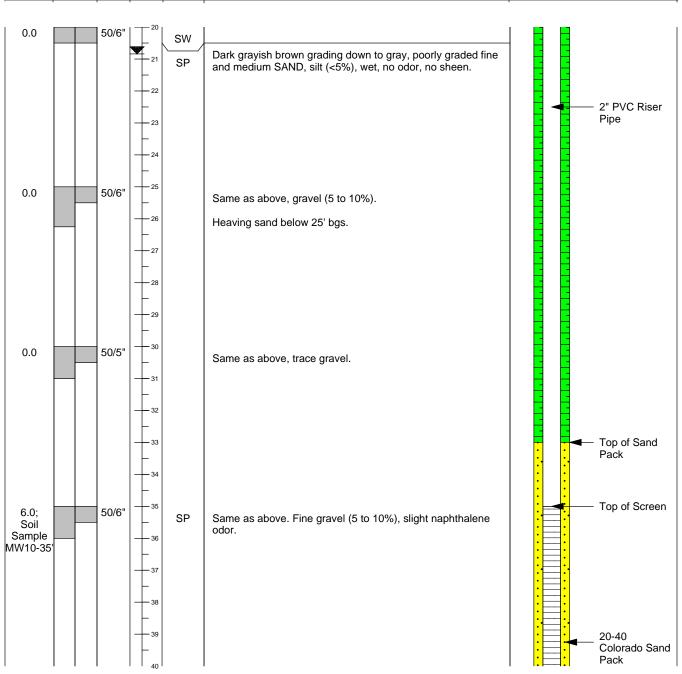
Drill Type: CME 75; 4-inch HSAClient: Bill SwensenSample Method: 18" D&M Split-spoonProject: Swensen-WCDBoring Diameter: 8"Site Location:3133 Cedar St,

Boring Depth (ft bgs): 46 ft bgs Groundwater ATD (ft bgs): 21.84*

Tacoma, WA.

Remarks: *Depth to water from top of casing from 6/22/10 groundwater sampling.

PID Reading	DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sample ID	RECOVERY	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



Ground Surf Elev. & Datum: 244.5 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 697945.61 Longitude/Easting: 1149387.71 Casing Elevation: 244.22 ft Monitoring Well ID: MW-10

Drill Date: June 10, 2010 **Logged By:** John LaManna

Drilled By: Curtis Askew / Cascade Drilling

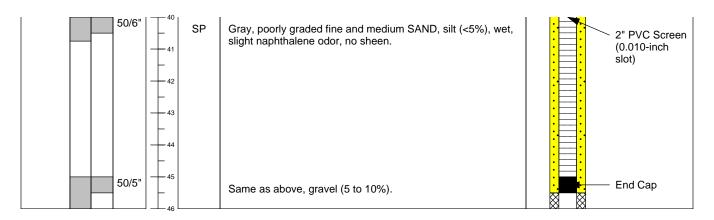
Drill Type: CME 75; 4-inch HSA
Sample Method: 18" D&M Split-spoon
Boring Diameter: 8"
Client: Bill Swensen
Project: Swensen-WCD
Site Location:3133 Cedar St,

Boring Depth (ft bgs): 46 ft bgs Groundwater ATD (ft bgs): 21.84*

Tacoma, WA.

Remarks: *Depth to water from top of casing from 6/22/10 groundwater sampling.

PID Reading	DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sample ID	RECOVERY	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



Ground Surf Elev. & Datum: 243.9 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 697807.34 Longitude/Easting: 1149380.41 Casing Elevation: 243.35 ft

Monitoring Well ID: MW-11

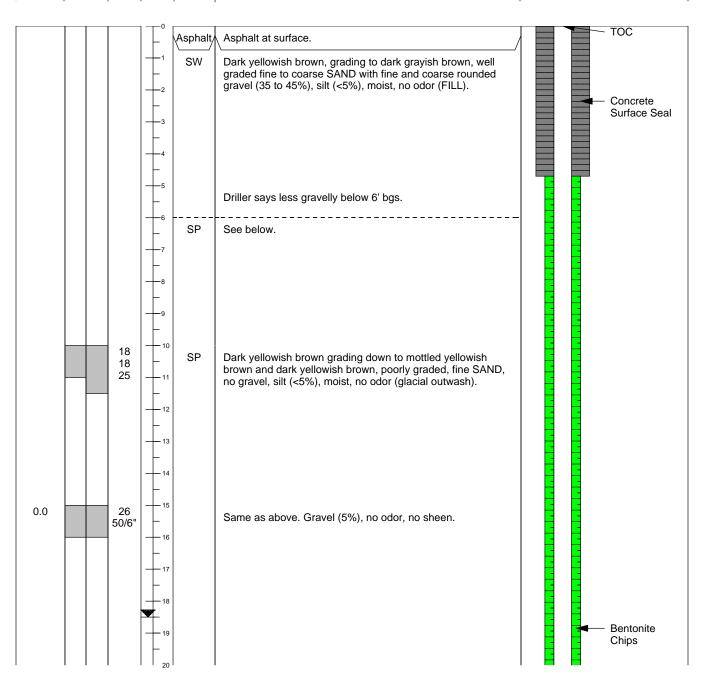
Drill Date: June 10, 2010 **Logged By:** John LaManna

Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA
Sample Method: 18" D&M Split-spoon
Boring Diameter: 8"
Client: Bill Swensen
Project: Swensen-WCD
Site Location:3133 Cedar St,

Boring Depth (ft bgs): 46 ft bgs Groundwater ATD (ft bgs): 18.5 Tacoma, WA.

PID	Reading	DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sa	imple ID	RECOVERY	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



Ground Surf Elev. & Datum: 243.9 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 697807.34 Longitude/Easting: 1149380.41 Casing Elevation: 243.35 ft

Monitoring Well ID: MW-11

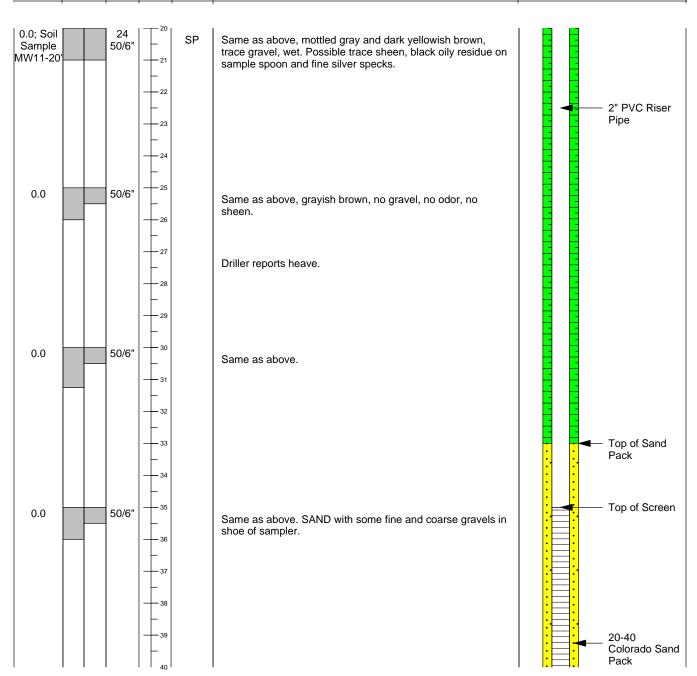
Drill Date: June 10, 2010 **Logged By:** John LaManna

Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA
Sample Method: 18" D&M Split-spoon
Boring Diameter: 8"
Client: Bill Swensen
Project: Swensen-WCD
Site Location:3133 Cedar St,

Boring Depth (ft bgs): 46 ft bgs Groundwater ATD (ft bgs): 18.5 Tacoma, WA.

	_					
PID Readin	g DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sample ID	RECOVERY	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



Ground Surf Elev. & Datum: 243.9 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 697807.34 Longitude/Easting: 1149380.41 Casing Elevation: 243.35 ft

Monitoring Well ID: MW-11

Drill Date: June 10, 2010 **Logged By:** John LaManna

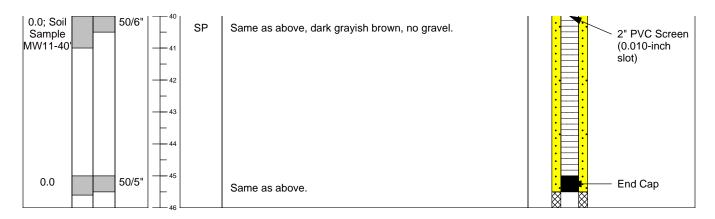
Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA
Sample Method: 18" D&M Split-spoon
Boring Diameter: 8"
Client: Bill Swensen
Project: Swensen-WCD
Site Location:3133 Cedar St,

Boring Depth (ft bgs): 46 ft bgs Groundwater ATD (ft bgs): 18.5

Tacoma, WA.

PID Readir	g DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sample II	RECOVER	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



Ground Surf Elev. & Datum: 244.5 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 698082.94 Longitude/Easting: 1149457.03 Casing Elevation: 243.97 ft Monitoring Well ID: MW-12

Drill Date: June 11, 2010 **Logged By:** John LaManna

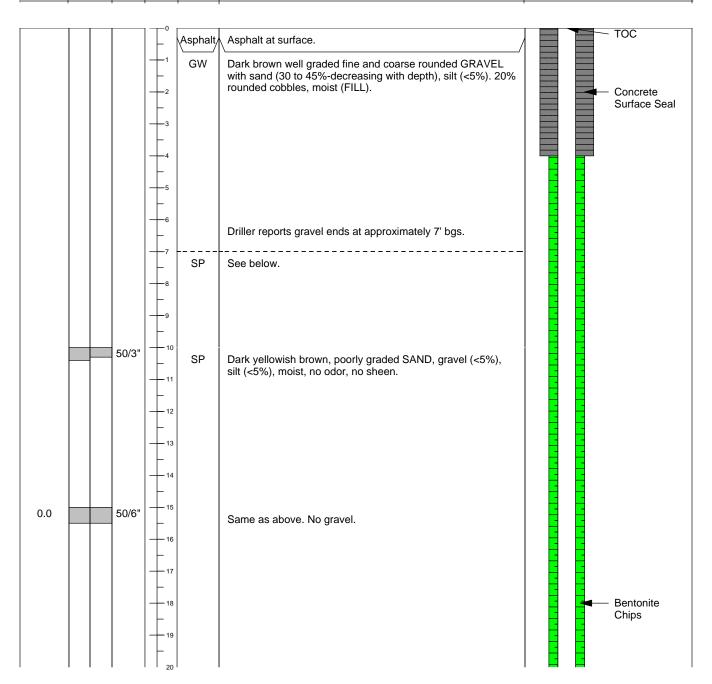
Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA
Sample Method: 18" D&M Split-spoon
Boring Diameter: 8"
Client: Bill Swensen
Project: Swensen-WCD
Site Location:3133 Cedar St,

Boring Depth (ft bgs): 47 ft bgs Groundwater ATD (ft bgs): 23.2

Tacoma, WA.

PID Reading	DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sample ID	RECOVERY	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



Ground Surf Elev. & Datum: 244.5 NGVD 29

Coordinate System: NAD83/98 Latitude/Northing: 698082.94 Longitude/Easting: 1149457.03 Casing Elevation: 243.97 ft

Monitoring Well ID: MW-12

Drill Date: June 11, 2010 **Logged By:** John LaManna

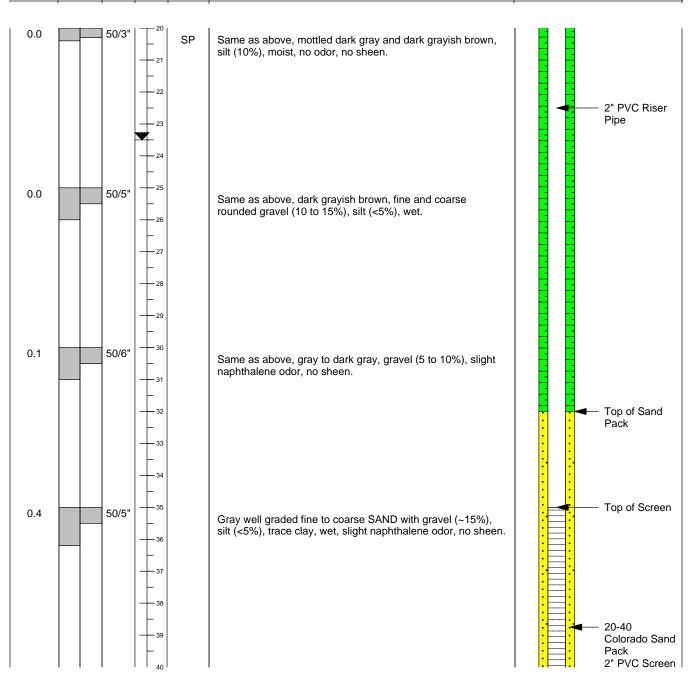
Drilled By: Curtis Askew / Cascade Drilling

Drill Type: CME 75; 4-inch HSA
Sample Method: 18" D&M Split-spoon
Boring Diameter: 8"
Client: Bill Swensen
Project: Swensen-WCD
Site Location:3133 Cedar St,

Boring Depth (ft bgs): 47 ft bgs Groundwater ATD (ft bgs): 23.2

Tacoma, WA.

PID Reading	DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sample ID	RECOVERY	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



Ground Surf Elev. & Datum: 244.5 NGVD 29

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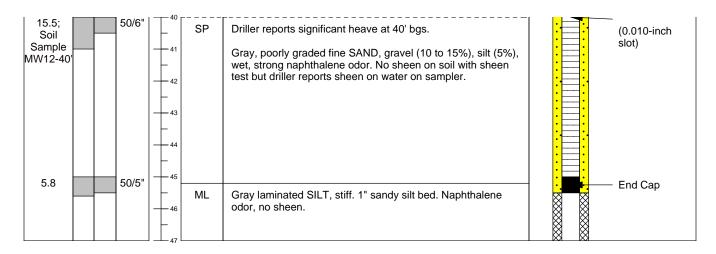
Drilled By: Curtis Askew / Cascade Drilling

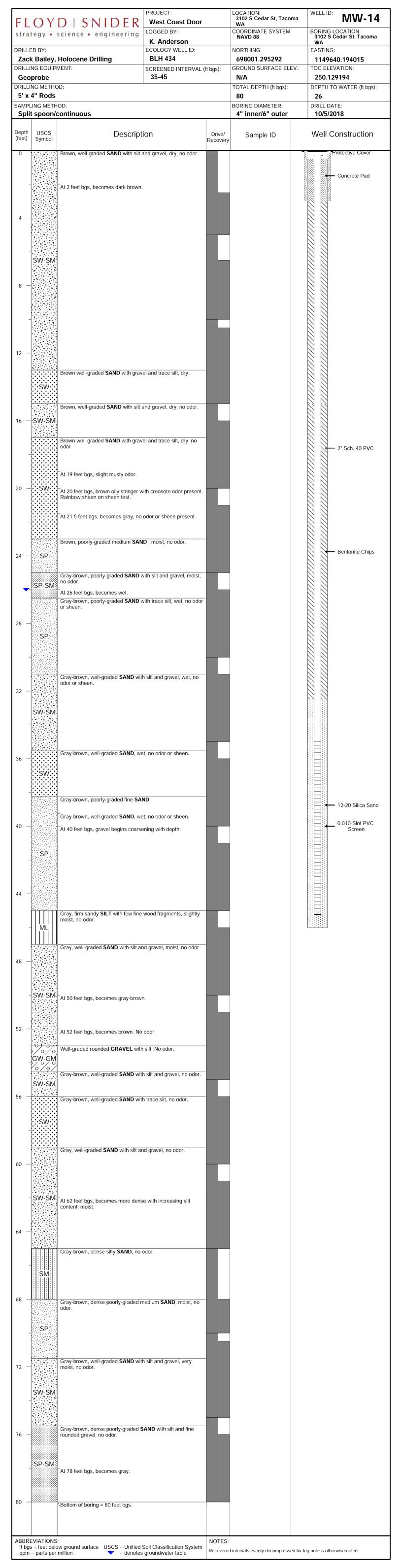
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Boring Depth (ft bgs): 47 ft bgs Groundwater ATD (ft bgs): 23.2

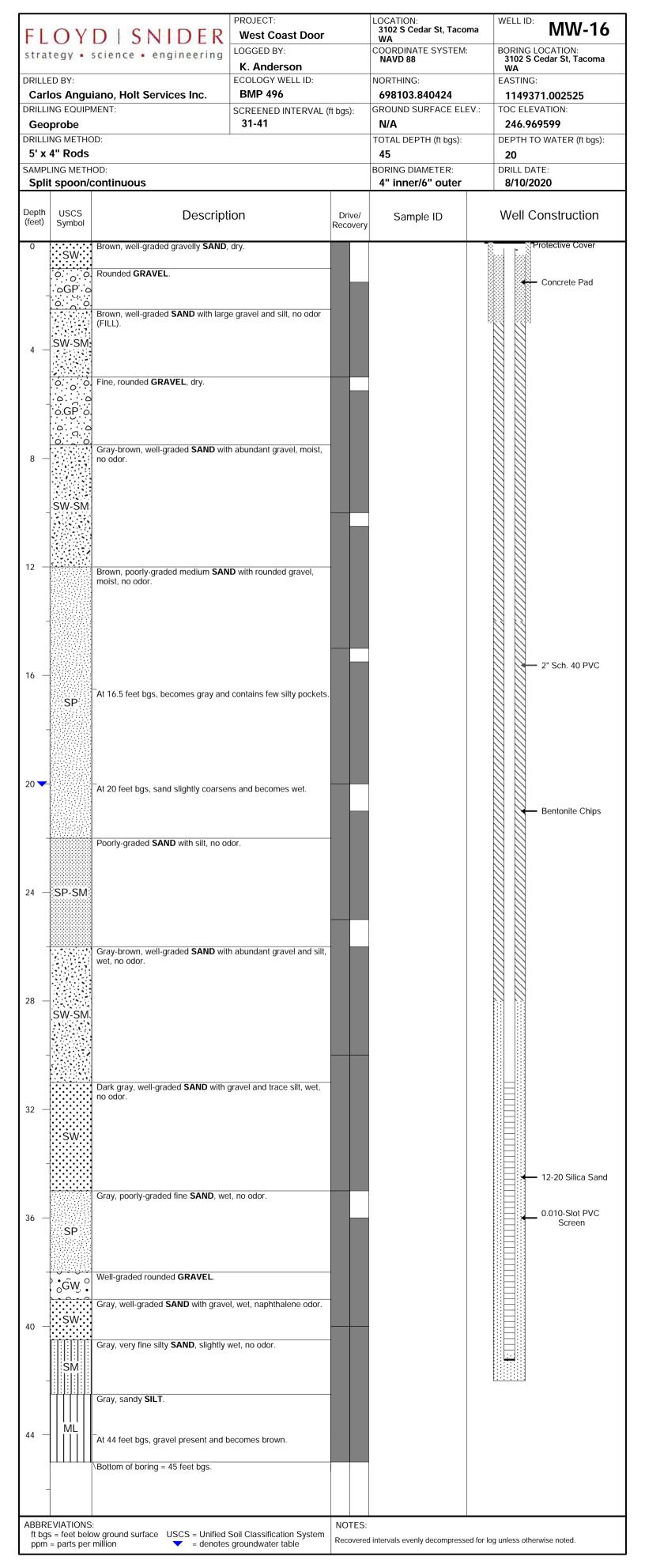
Tacoma, WA.

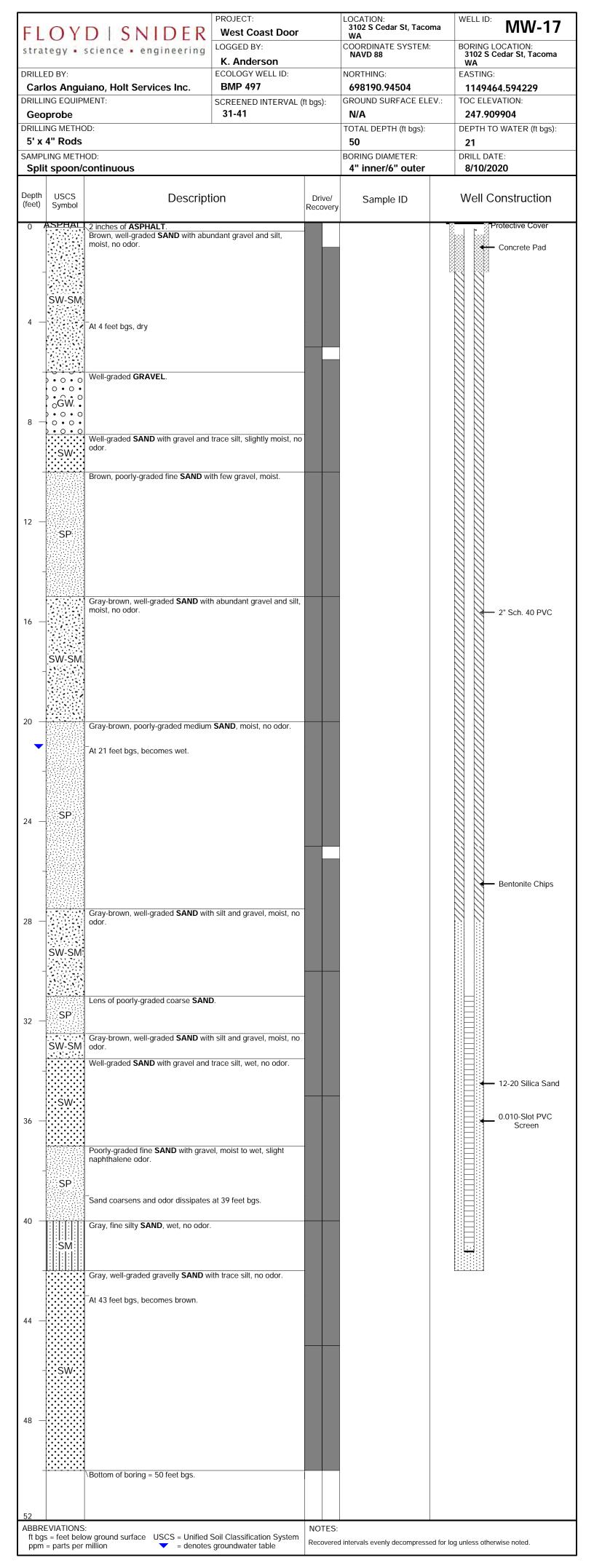
PID Readir	g DRIVE /	BLOW	DEPTH	USCS	SOIL DESCRIPTION AND OBSERVATIONS: (color, texture,	MONITORING WELL
/Sample II	RECOVER	COUNT	FT BGS	SYMBOL	moisture, MAJOR CONSTITUENT, odor, staining, sheen, debris, etc.)	DETAIL



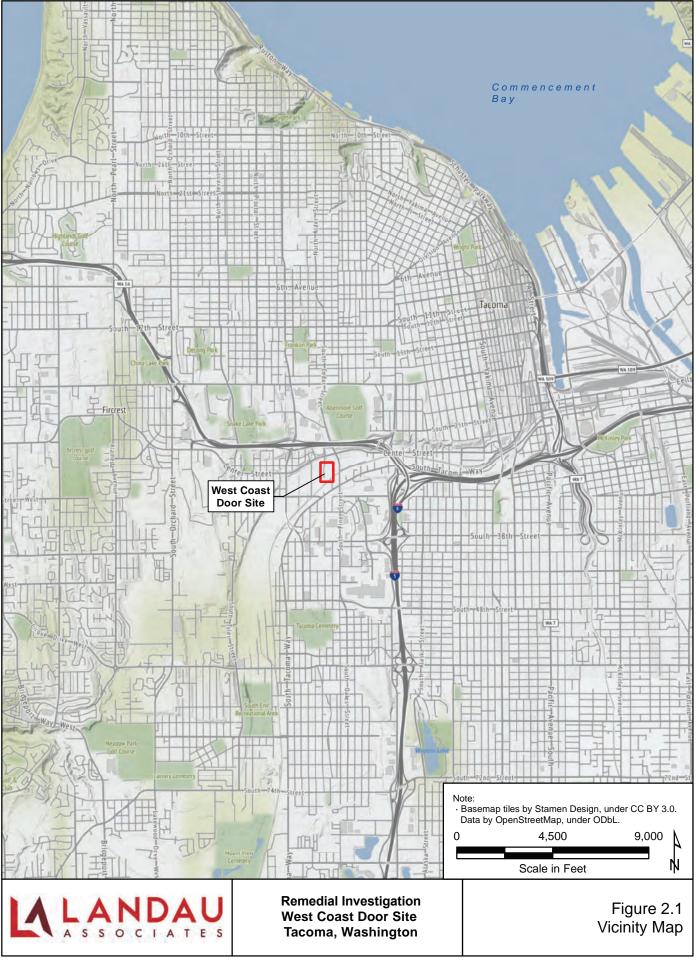


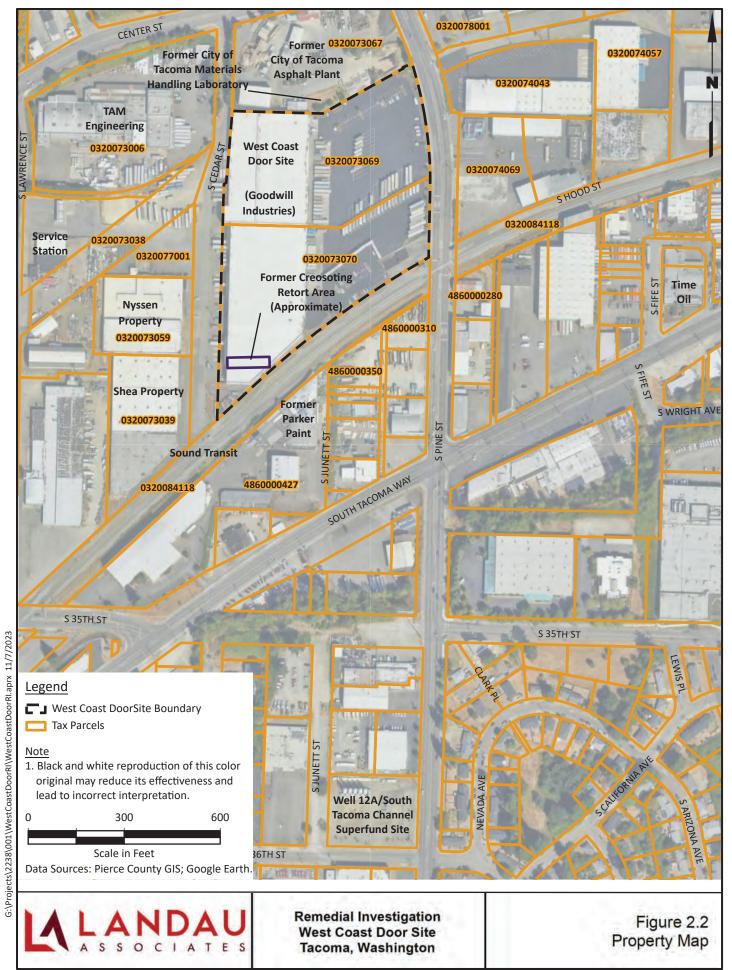
ΕT	OV	DISNIDER	PROJECT: West Coast Door		LOCATION: 3102 S Cedar St, Tacol WA	well id:	15
		science • engineering	LOGGED BY: K. Anderson		COORDINATE SYSTEM: NAVD 88	BORING LOCATION: 3102 S Cedar St, Taco WA	oma
		n, Holt Services, Inc.	ECOLOGY WELL ID: BMP 498	5. J	NORTHING: 697918.299897	EASTING: 1149273.549423	
Geo	NG EQUIPI probe NG METHO		SCREENED INTERVAL (1 33.5-43.5	it bgs):	GROUND SURFACE ELE N/A TOTAL DEPTH (ft bgs):	EV.: TOC ELEVATION: 245.277262 DEPTH TO WATER (ft b.	us).
5' x	4" Rods ING METH				50 BORING DIAMETER:	20 DRILL DATE:	
	-	continuous			4" inner/6" outer	8/11/2020	
Depth (feet)	USCS Symbol	Descript	ion	Drive/ Recovery	Sample ID	Well Constructio	
0		Brown, well-graded SAND with well and silt, no odor.	-graded rounded gravel			Concrete Pac	
-							
4 —							
	SW-SM						
_							
8 —		At 7 feet bgs, becomes moist.					
0		Brown, poorly-graded medium SAN	ID with silt and gravel, no				
_	SP-SM	At 10 to 15 feet bgs, sample not rec	covered.				
12 —							
12							
_							
16 —		Brown, poorly-graded medium SAN odor.	ID with silt and gravel, no			Bentonite Chi	ips
10		At 17 feet bgs, few peaty pockets.				2" Sch. 40 PV	/C
_	SP-SM						
20 🕶							
20		At 20 feet bgs, becomes wet. Poorly-graded medium SAND with	trace silt, wet, no odor.				
=		At 22 feet bgs, large gravel present					
24							
24 —							
-	SP	At 26 feet bgs, gravel content decre	eases with little gravel				
28 —							
20							
_						Bentonite Chi	ips
32 —		Brown, well-graded SAND with grav	vel, wet, no odor.				
JZ —	SW						
-	∵ GW	Lens of well-graded gravel Brown, well-graded SAND with grav	vel, wet, no odor.				
36 —	. SW.	_					
		Brown, poorly-graded medium SAN	ID, wet, no odor.				
-						12-20 Silica S 0.010-Slot PV Screen	
40 —							
	SP						
-							
44 —		Gray, well-graded very fine SAND , naphthalene odor.					
	::SW:::	Gray, well-graded SAND with silt an odor. Gray, poorly-graded fine SAND , po coarsens with depth.					
-		обавона міш исриі.					
48 —	SP						
	C/M-	Gray, well-graded SAND with silt an	nd gravel, no odor, wet.				
_	SW	\Bottom of boring = 50 feet bgs.					
52							
ft bgs	EVIATIONS = feet belo = parts per	ow ground surface USCS = Unified	Soil Classification System groundwater table	NOTES: Recovered	l intervals evenly decompresse	ed for log unless otherwise noted.	





Selected Remedial Investigation Figures





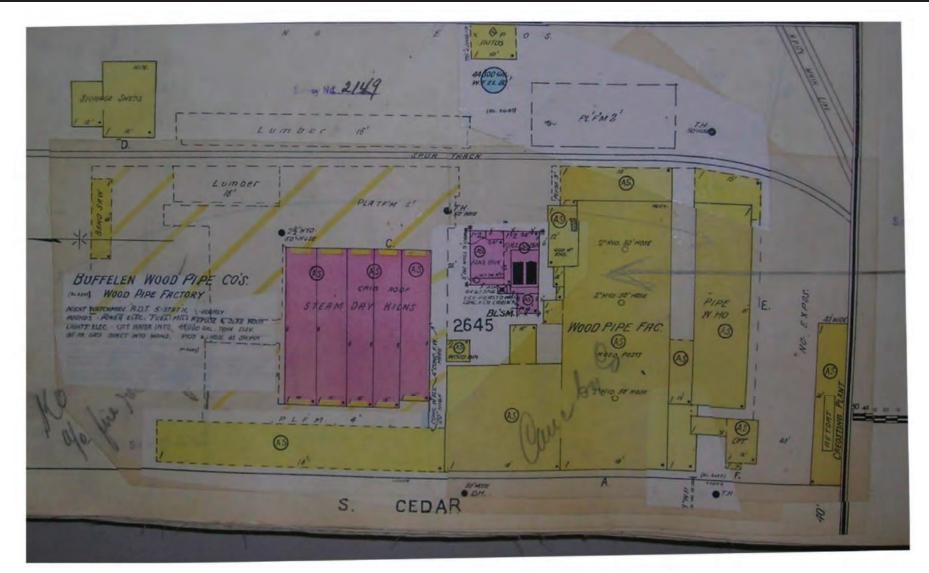
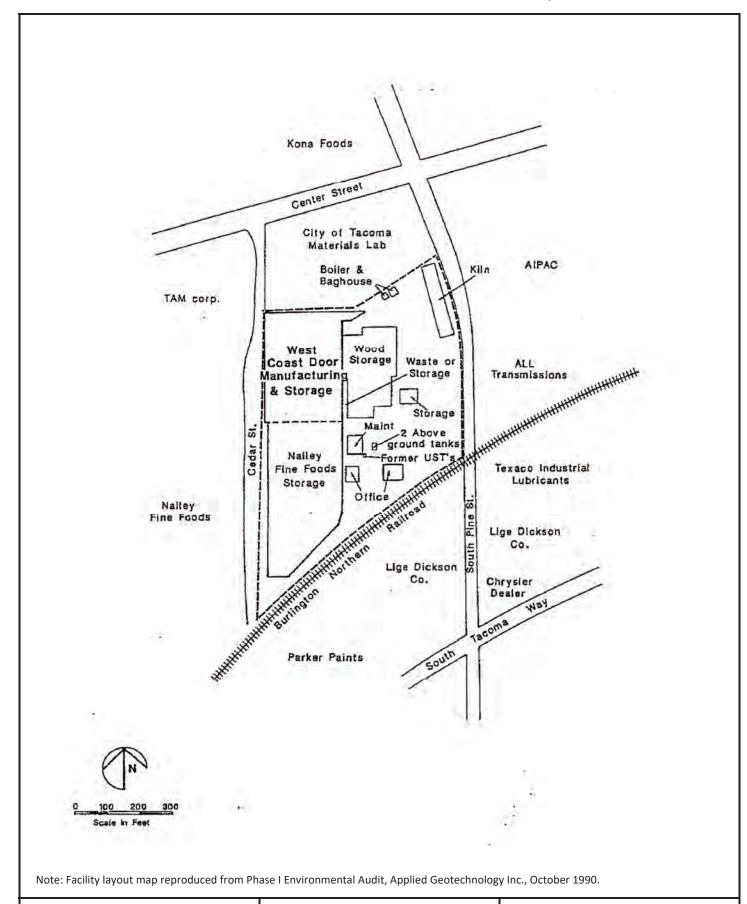


Image from Sanborn Maps, 1912



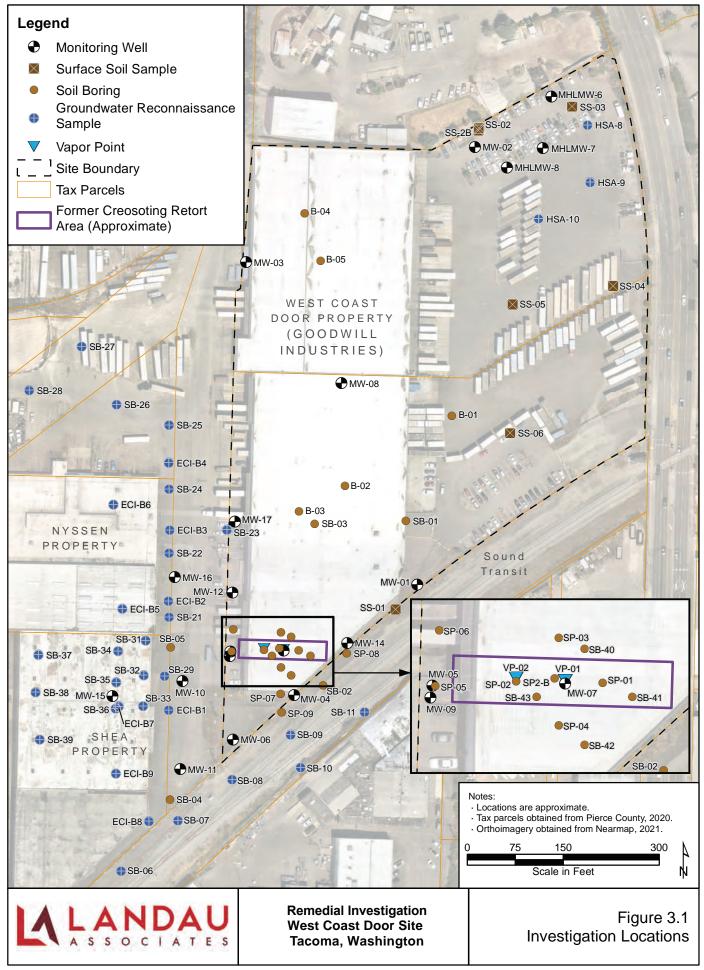
Remedial Investigation West Coast Door Site Tacoma, Washington

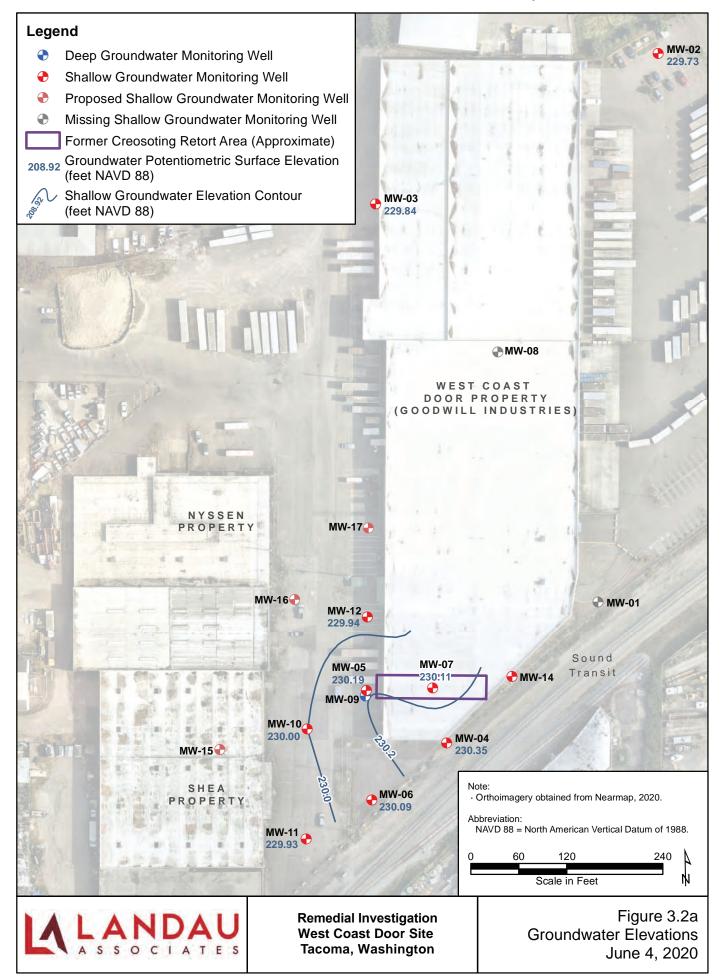
Figure 2.3 Former Creosoting Facility Operations

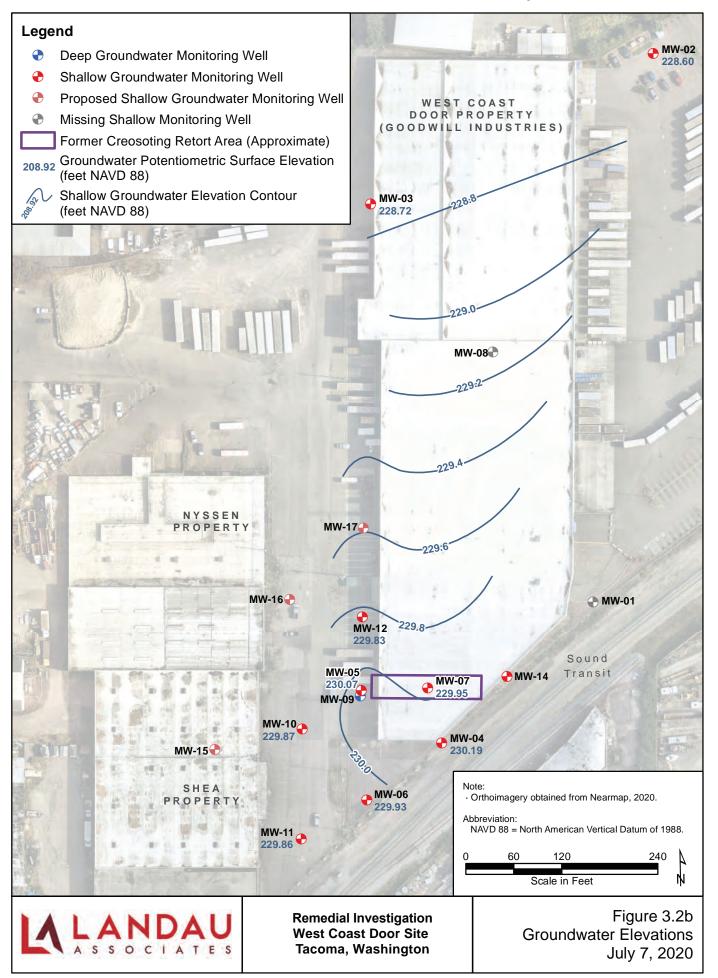


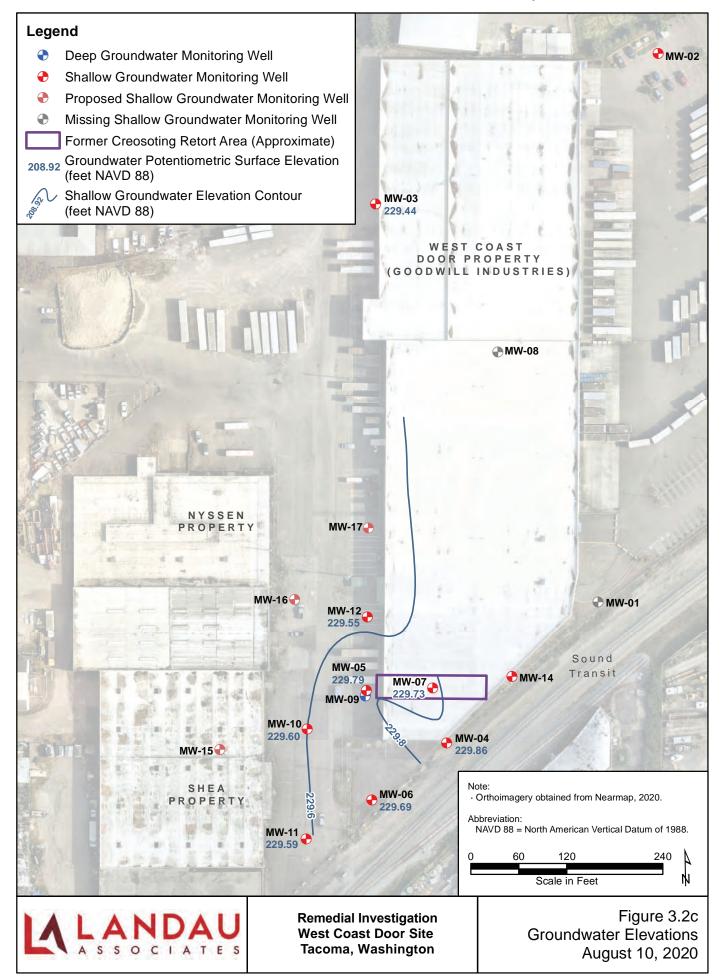


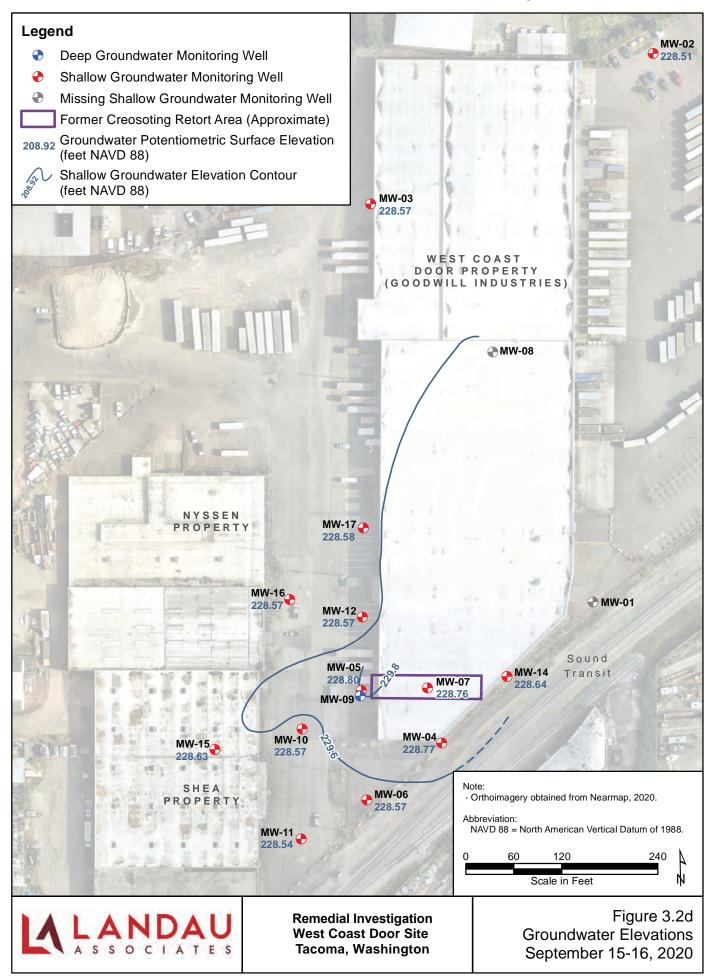
Remedial Investigation West Coast Door Site Tacoma, Washington Figure 2.4 Former West Coast Door Facility Operations

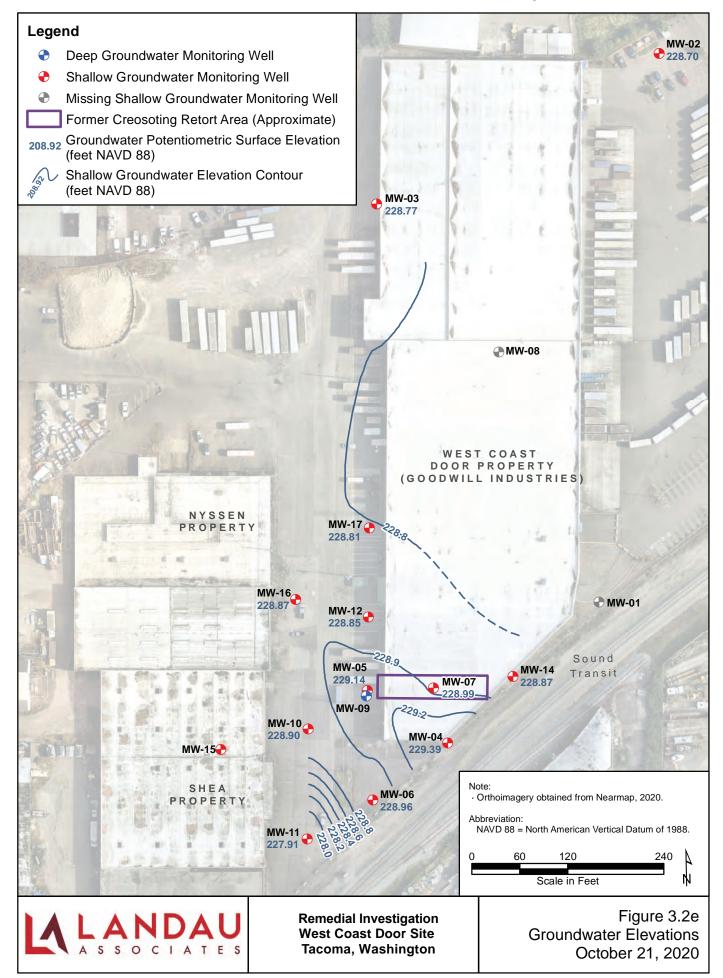


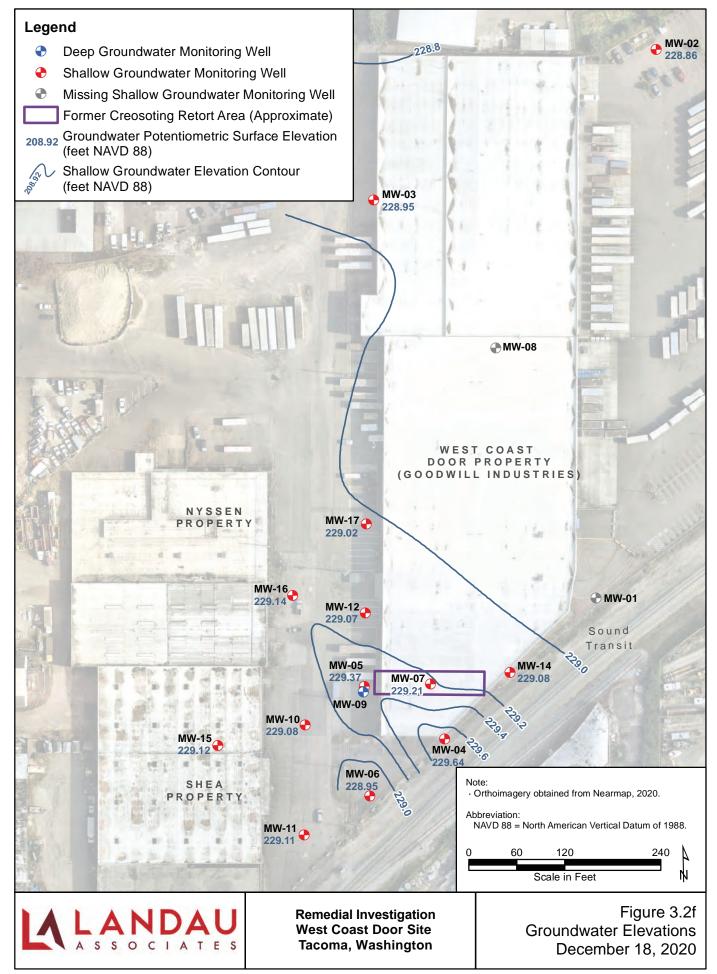


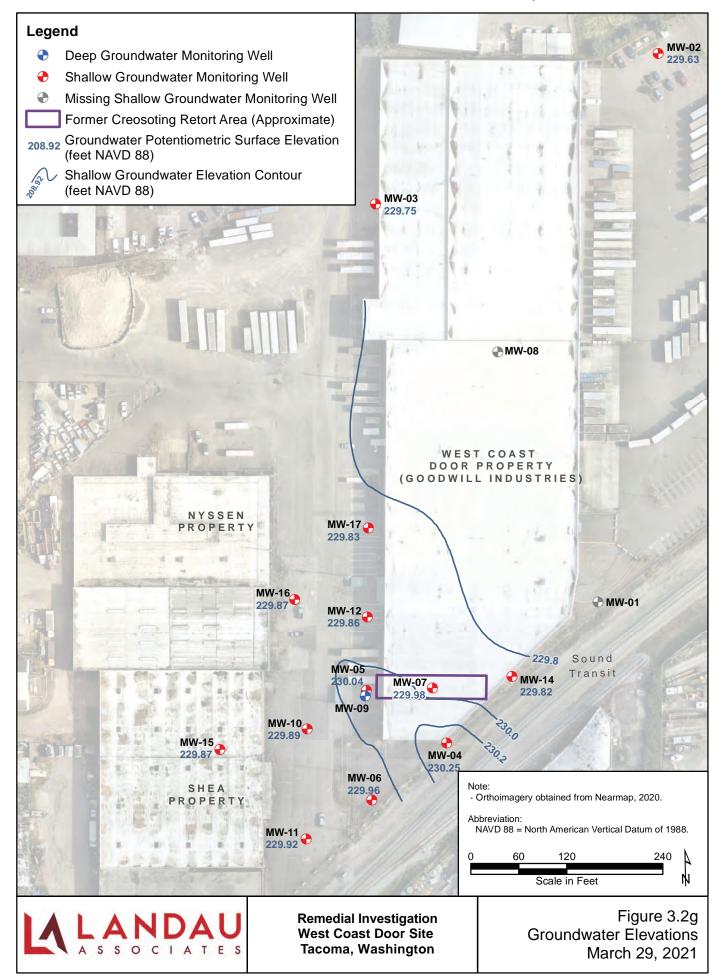


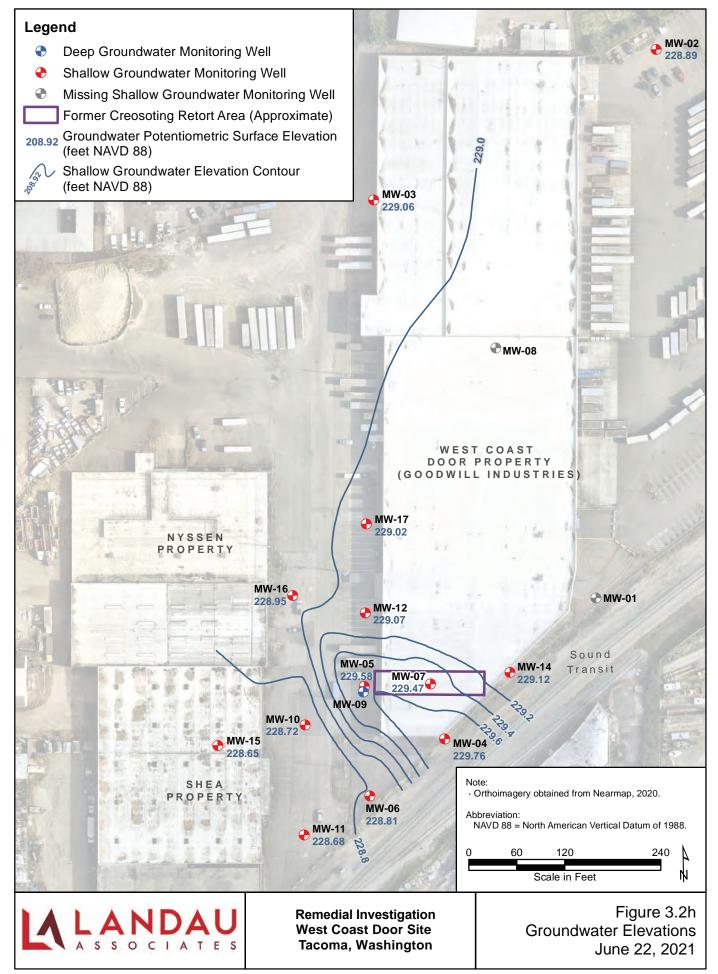












Legend **Location Type** Shallow Groundwater Monitoring Well Groundwater Reconnaissance Sample⁽¹⁾ Conc. in µg/L (Dashed Where Inferred) Total Naphthalene >160⁽²⁾-≤1,600 µg/L Total Naphthalene >1,600-≤16,000 µg/L Total Naphthalene >16,000 µg/L **Other Features** Site Boundary Tax Parcels Former Creosoting Retort Area (Approximate) WEST COAST DOOR PROPERTY (GOODWILL INDUSTRIES) MW-17 70 NYSSEN Transit PROPERTY MW-16 1,600 MW-12 1. Groundwater reconnaissance sample results, collected in 14,000 April and May, 2009, are shown only where they delineate the extent of detectable contamination in areas inaccessible for installation of permanent monitoring wells. The proposed Site cleanup level for naphthalene is MW-14 160 µg/L. Results in **bold** exceed this level. 3,600 The extent of the >16,000 μ g/L is inferred in the vicinity of the Former Creosoting Retort Area because monitoring wells are not screened with sufficient depth MW-10 🔴 to sample the most highly contaminated groundwater 17,000 MW-15 interval SB-11 1.0 U Only wells screened in the most highly contaminated groundwater interval of 35-45 ft bgs were sampled. 1.0 U SB-9 All results are in µg/L. MW-06 SHEA Tax parcels obtained from Pierce County, 2020. 4,500 3.800 PROPERTY Orthoimagery obtained from Nearmap, 2021. \$B-10 MW-11 Abbreviation: \$B-08 1.0 U 1.0 U 15.0 bgs = Below ground surface μg/L = Micrograms per liter \$B-07 1.0 U U = Analyte was not detected at the given reporting limit 75 \$B-06 1.3 Scale in Feet Figure 5.1

LANDAU ASSOCIATES Remedial Investigation West Coast Door Site Tacoma, Washington Figure 5.1 Naphthalene in Shallow Groundwater (35-45 ft bgs), December 2020

Modified from Floyd | Snider 2022 draft RI/FS.

Legend **Location Type** Shallow Groundwater Monitoring Well Groundwater Reconnaissance Sample⁽¹⁾ Conc. in µg/L (Dashed Where Inferred) Total Naphthalene >160⁽²⁾-≤1,600 µg/L Total Naphthalene >1,600-≤16,000 µg/L Total Naphthalene >16,000 µg/L **Other Features** Site Boundary Tax Parcels Former Creosoting Retort Area (Approximate) WEST COAST DOOR PROPERTY (GOODWILL INDUSTRIES) MW-17 Sound 1.8 NYSSEN Transit PROPERTY MW-16 Notes: 75 Groundwater reconnaissance sample results, collected in April and May, 2009, are shown only where they delineate the extent of detectable contamination in areas MW-12 inaccessible for installation of permanent monitoring wells 17,000 MW-14 The proposed Site cleanup level for nahpthalene is 160 μ g/L. Results in **bold** exceed this level. The extent of the >16,000 μ g/L is inferred in the vicinity 7.900 of the Former Creosoting Retort Area because monitoring wells are not screened with sufficient depth to sample the most highly contaminated groundwater MW-10 MW-15 interval **SB-11** 1.5 16,000 Only wells screened in the most highly contaminated groundwater interval of 35-45 ft bgs were sampled. 1.0 U SB-09 All results are in µg/L. MW-06 SHEA Tax parcels obtained from Pierce County, 2020. 4,500 6,500 PROPERTY Orthoimagery obtained from Nearmap, 2021. \$B-10 MW-11 🕀 Abbreviation: \$B-08 1.0 U 1.0 U bgs = Below ground surface 1.3 U μg/L = Micrograms per liter \$B-07 1.0 U U = Analyte was not detected at the given reporting limit 300 75 \$B-06 Scale in Feet

LANDAU ASSOCIATES Remedial Investigation West Coast Door Site Tacoma, Washington Figure 5.2 Naphthalene in Shallow Groundwater, (35-45 ft bgs), June 2021

Legend **Location Type** Shallow Groundwater Monitoring Well Groundwater Reconnaissance Sample(1) Conc. in µg/L (Dashed Where Inferred) Benzene >5 and Total Xylenes >1,000(2) Other Features ¹ Site Boundary Tax Parcels Former Creosoting Retort Area (Approximate) WEST COAST DOOR PROPERTY (GOODWILL INDUSTRIES) Sound MW-17 Transit B:1.0 U NYSSEN X:1.0 U PROPERTY 1. Groundwater reconnaissance sample results, collected in May 2009, are shown only where they MW-16 MW-12 delineate the extent of detectable contamination in B:2.7 areas inaccessible for installation of permanent wells. B:9.3 X:260 2. The proposed Site cleanup level for benzene is X:2.000 5 μg/L and for total xylenes is 1,000 μg/L. Results in bold exceed this level. MW-14 The extent of the >5 μ g/L for benzene and >1,000 μ g/L for total xylenes is inferred in the vicinity of the Former B:1.0 U Creosoting Retort Area because monitoring wells are X:68.0 not screened with sufficient depth to sample the most highly contaminated groundwater interval. MW-10 MW-15 Only wells screened in the most highly contaminated B:10.0 U groundwater interval of 35-45 ft bgs were sampled. All results are in $\mu g/L$. B:1.0 U X:1,300 SB-11 B: 0.20 U X:1.0 U Tax parcels obtained from Pierce County, 2020. **SB-09** MW406 X: 0.40 U Orthoimagery obtained from Nearmap, 2021. SHEA **⊕** B: **20 U** B:1.0 U PROPERTY X: 40 U X:18.0 Abbreviations: SB-10 MW-11 B = Benzene B: 1.0 U SB-08 B: 0.20 U B: 0.72 bgs = Below ground surface X:1.0 U X: 0.40 U μg/L = Micrograms per liter X: 0.40 U X = Total Xylenes SB-07 **⊕** B: 0.20 U X: 0.40 U U = Analyte was not detected at the given reporting limit SB-06 75 150 300 B: 0.20 U X: 0.40 U Scale in Feet Figure 5.3 **Remedial Investigation**

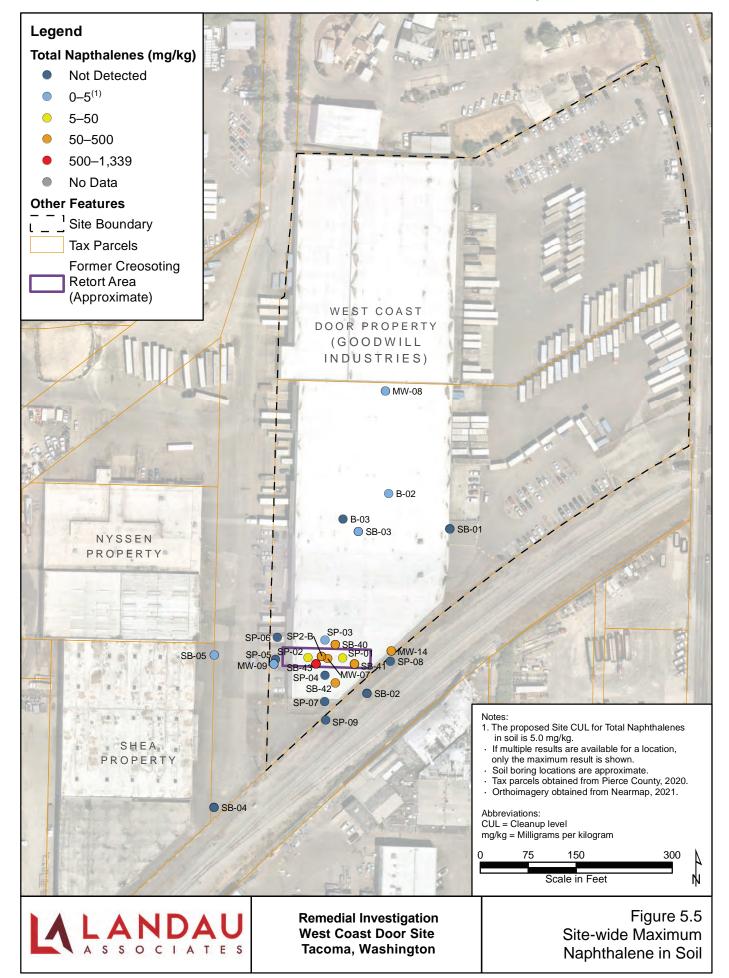
LANDAU ASSOCIATES Remedial Investigation
West Coast Door Site
Tacoma, Washington

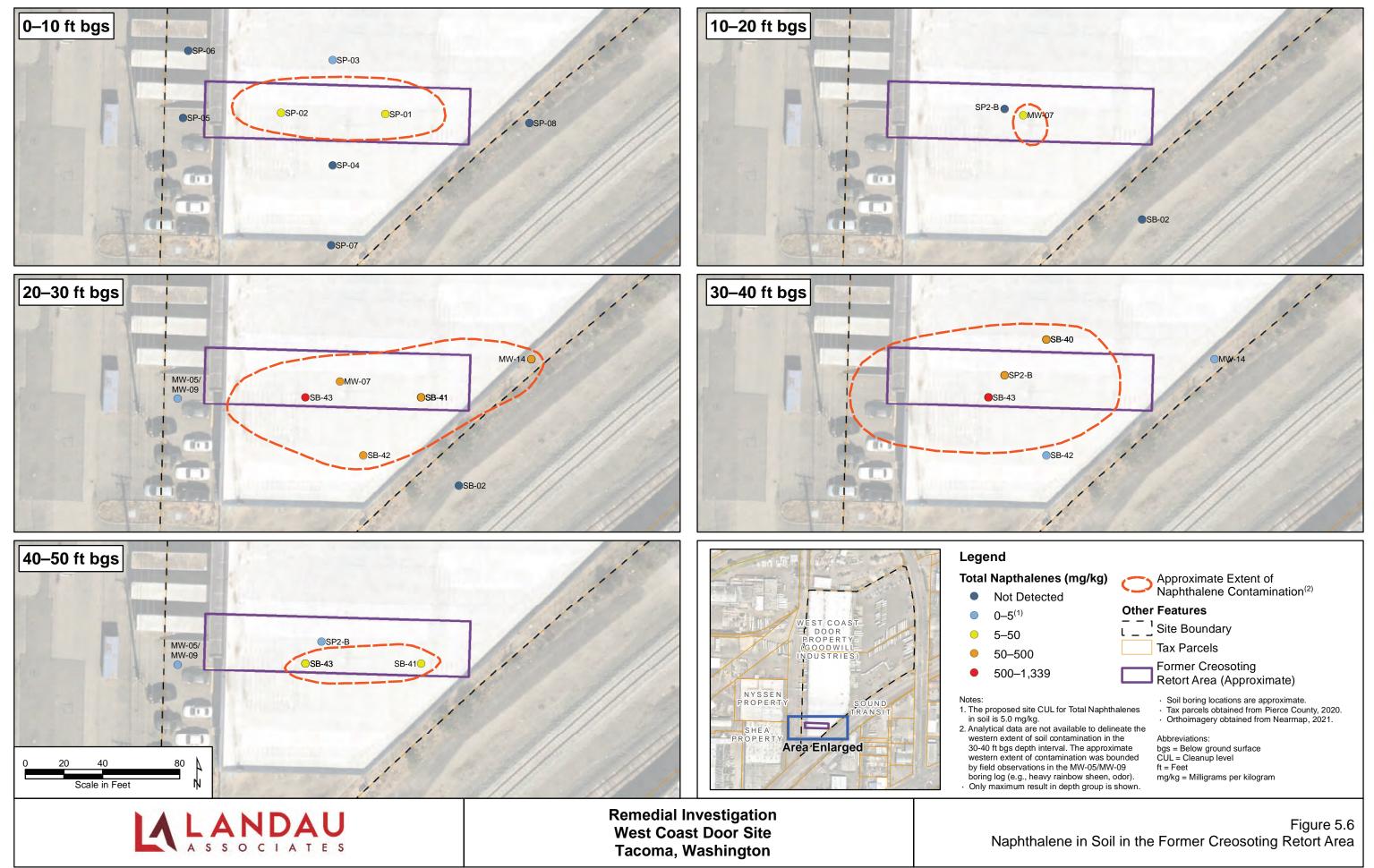
Figure 5.3 Benzene and Xylenes in Shallow Groundwater (35-45 ft bgs), December 2020

Legend **Location Type** Shallow Groundwater Monitoring Well Groundwater Reconnaissance Sample⁽¹⁾ Conc. in µg/L (Dashed Where Inferred) Benzene >5 and Total Xylenes >1,000⁽²⁾ Other Features ¹ Site Boundary Tax Parcels Former Creosoting Retort Area (Approximate) WEST COAST DOOR PROPERTY (GOODWILL INDUSTRIES) MW-17 Transit B:0.44 U NYSSEN X:1.0 U PROPERTY 1. Groundwater reconnaissance sample results, MW-16 collected in May 2009, are shown only where they B:0.44 U MW-12 delineate the extent of detectable contamination in areas inaccessible for installation of permanent wells. X:1.0 U B:11.0 2. The proposed Site cleanup level for benzene is X:10.0 U 5 μg/L and for total xylenes is 1,000 μg/L. Results in bold exceed this level. MW-14 The extent of the >5 μ g/L for benzene and >1,000 μ g/L for total xylenes is inferred in the vicinity of the Former B:4.4 U Creosoting Retort Area because monitoring wells are X:10.0 U not screened with sufficient depth to sample the most MW-10 highly contaminated groundwater interval. B:**8.8 U** Only wells screened in the most highly contaminated MW-15 SB-11 groundwater interval of 35-45 ft bgs were sampled. All results are in $\mu g/L$. X:20.0 U B:0.44 U B: 0.20 U X:1.0 U Tax parcels obtained from Pierce County, 2020. SB-09 X: 0.40 U MW-06 B: 20 U Orthoimagery obtained from Nearmap, 2021. SHEA B:4.4 U X: 40 U PROPERTY X:10.0 U Abbreviations: SB-10 MW-11 B:0.44 U B = Benzene **SB-08** B: 0.72 bgs = Below ground surface **⊕** B: 0.20 U X: 0.40 U X:1.0 U X: 0.40 U μg/L = Micrograms per liter X = Total Xylenes **SB-07** ⊕ B: 0.20 U X: 0.40 U U = Analyte was not detected at the given reporting limit **SB-06** 75 150 300 B: 0.20 U X: 0.40 U Scale in Feet Figure 5.4 **Remedial Investigation**

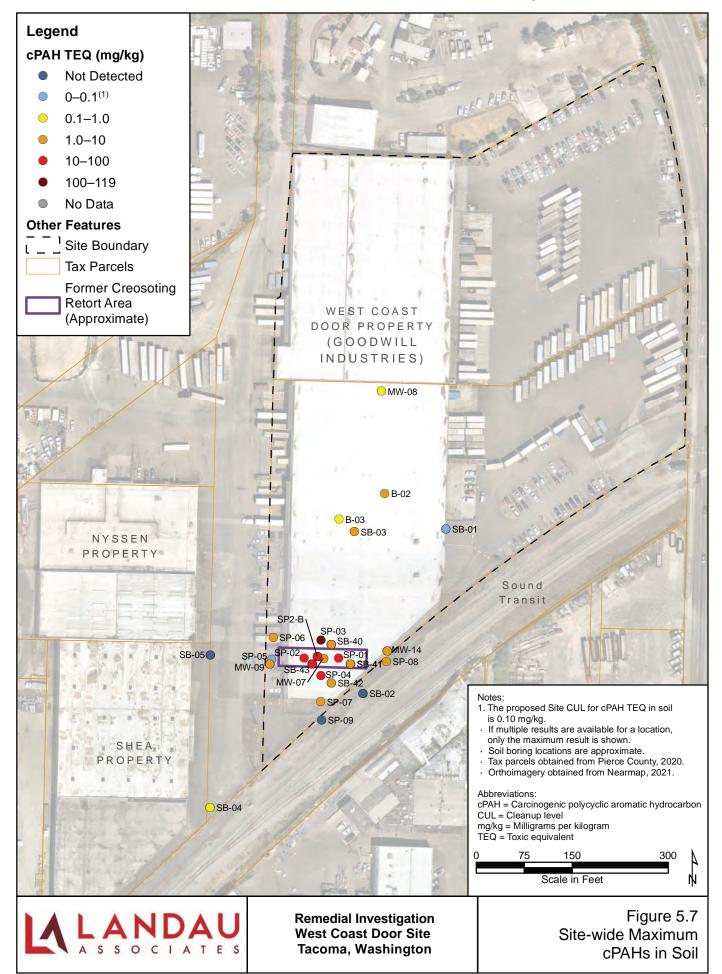
LANDAU ASSOCIATES Remedial Investigation
West Coast Door Site
Tacoma, Washington

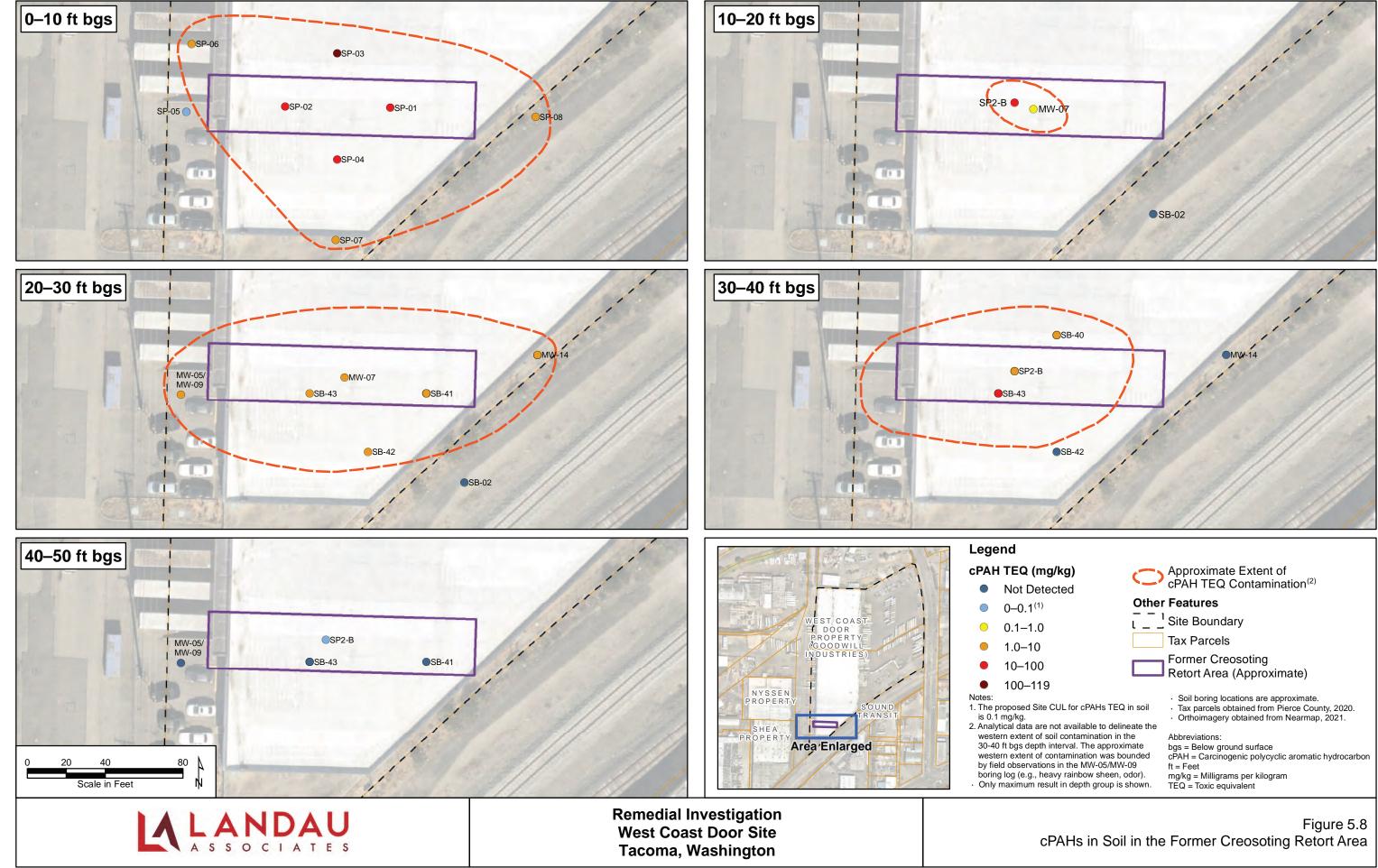
Figure 5.4 Benzene and Xylenes in Shallow Groundwater (35-45 ft bgs), June 2021



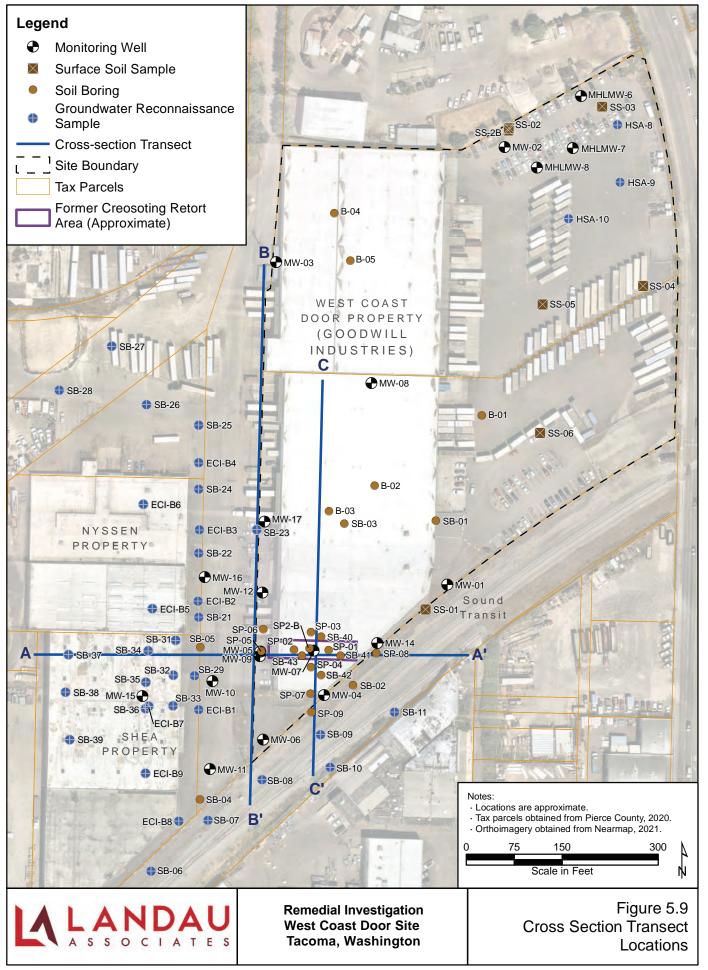


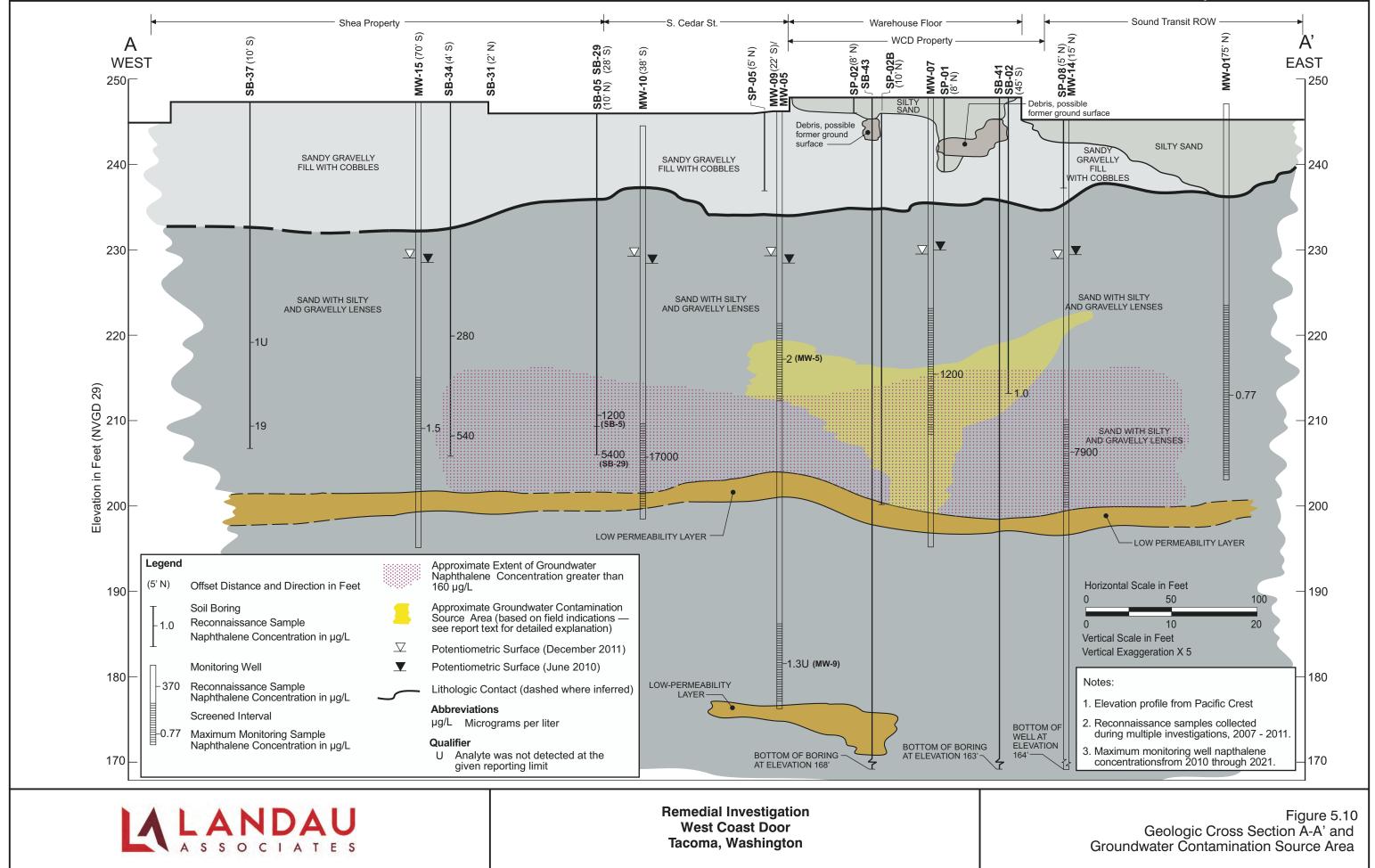
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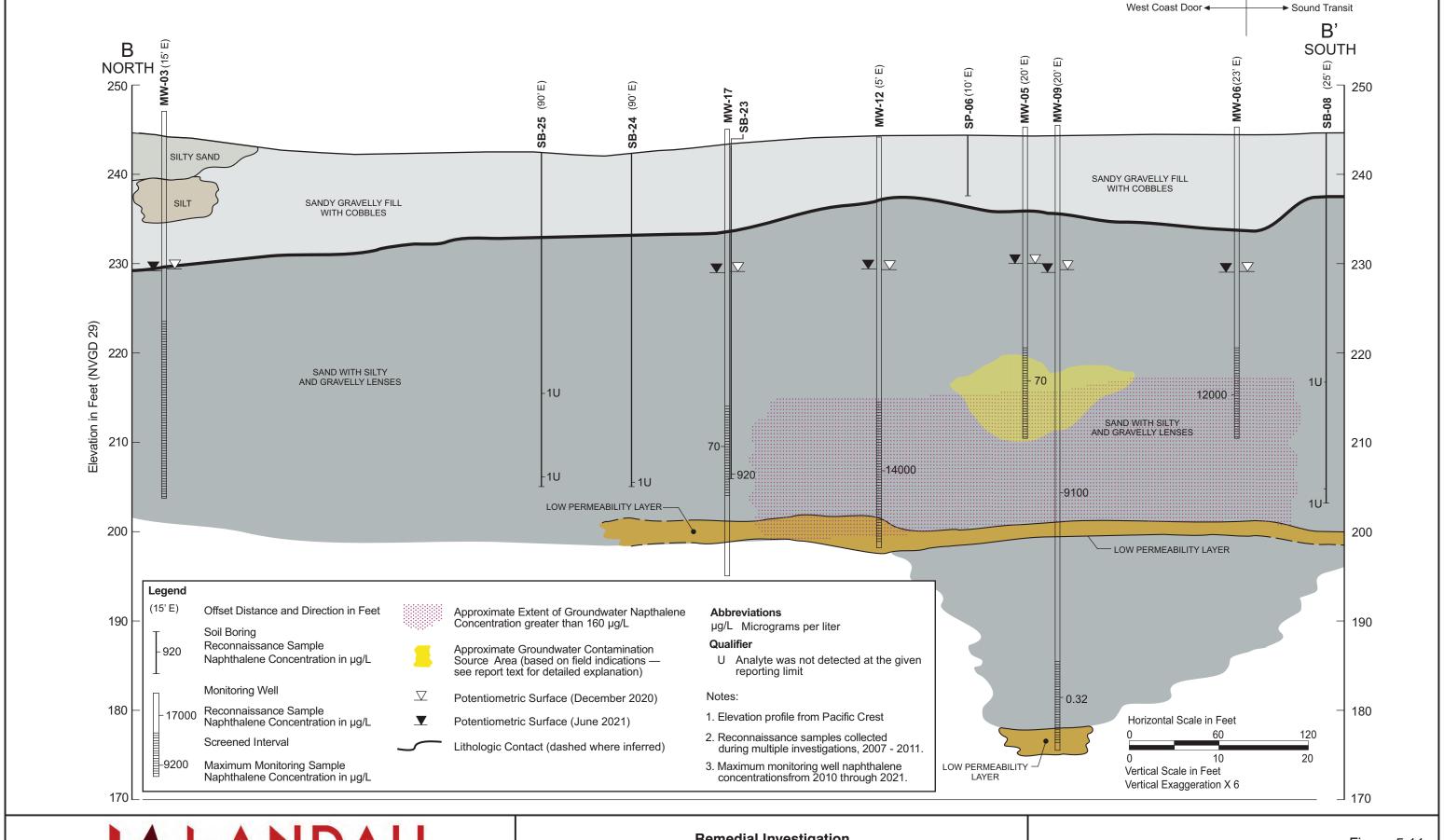




11/8/2023 P:\2238\001\R\Remedial Investigation\Figures



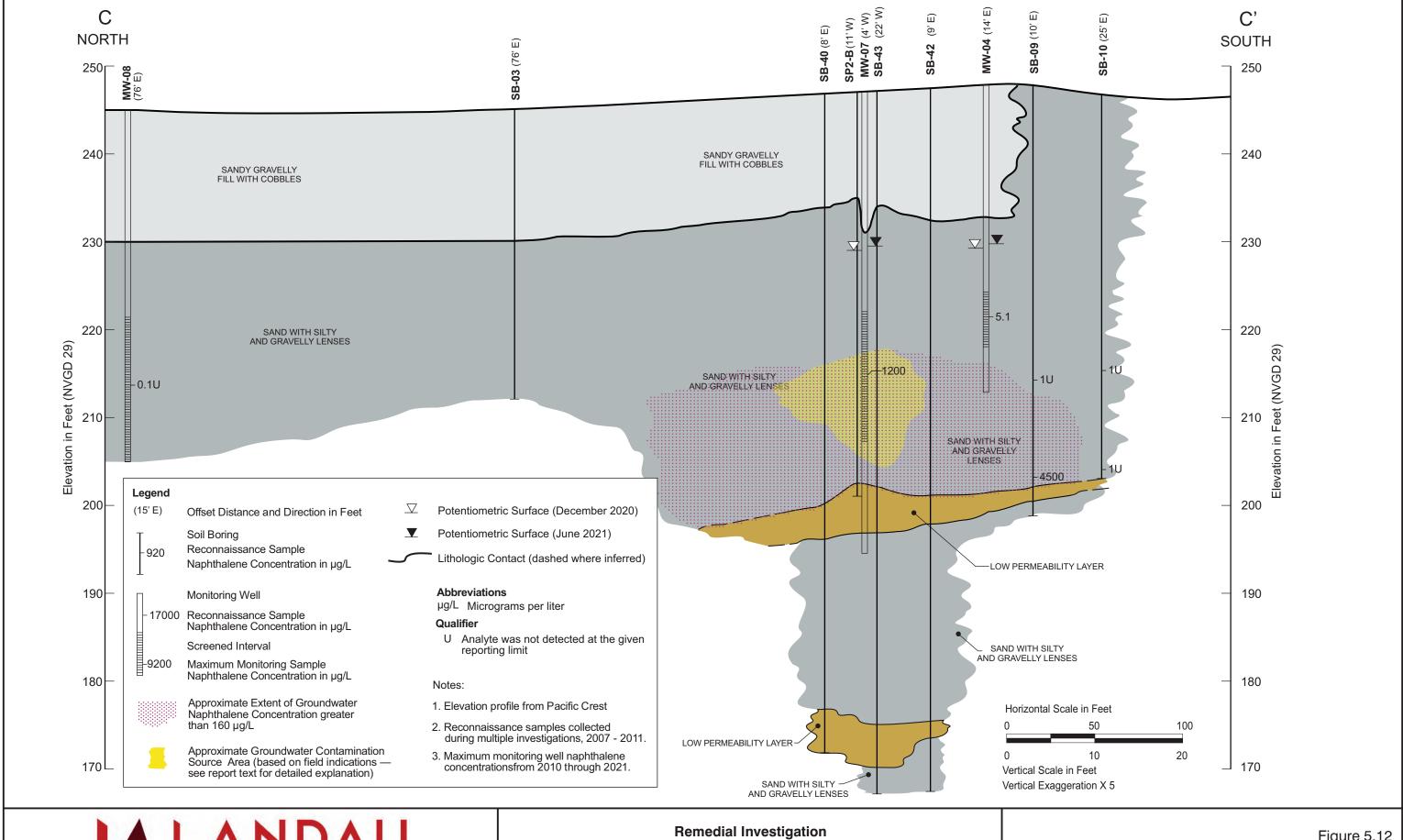




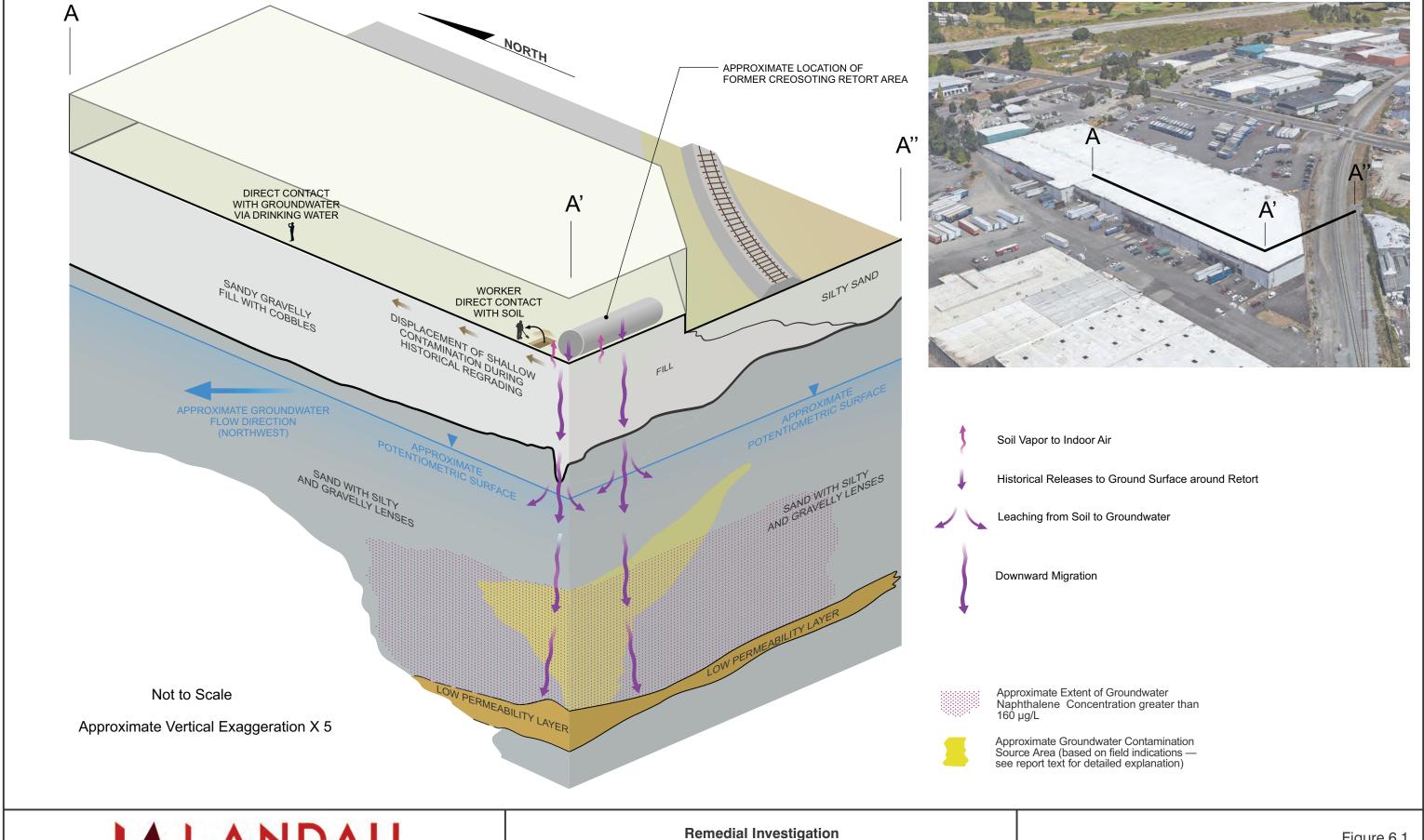
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Remedial Investigation West Coast Door Tacoma, Washington

Figure 5.11 Geologic Cross Section B-B' and Groundwater Contamination Source Area



LANDAU ASSOCIATES Remedial Investigation West Coast Door Tacoma, Washington Figure 5.12 Geologic Cross Section C-C' and Groundwater Contamination Source Area



LANDAU

Remedial Investigation West Coast Door Tacoma, Washington Figure 6.1 Conceptual Site Model

Selected Remedial Investigation Tables



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		2.34	MW-01	. ,	MW-02 MW-04				
			11100 11111 20								
		Sample Name	WCD-MW1-38	WCD-MW01-112712	WCD-MW01-030513	WCD-MW01-060413	WCD-MW01-100113	MW02-WL-03292021	WCD-MW4-33	WCD-MW04-120511	MW04-WL-03292021
		Sample Date	6/22/2010	11/27/2012	3/5/2013	6/4/2013	10/1/2013	3/29/2021	6/21/2010	12/5/2011	3/29/2021
Analyte	CAS No.	Screening Level (1)									
Semivolatile Organic Compounds 1-Methylnaphthalene	90-12-0	1		0.10 U				l	1		
2,4,5-Trichlorophenol	95-95-4	1,600		0.10 0							
	88-06-2	8.0									
2,4,6-Trichlorophenol	120-83-2	48							10 U		
2,4-Dichlorophenol									1		
2,4-Dimethylphenol	105-67-9	320							10 U		
2,4-Dinitrophenol	51-28-5	32							30 U		
2-Chloronaphthalene	91-58-7								40.11		
2-Chlorophenol	95-57-8	40							10 U		
2-Methylnaphthalene	91-57-6	-		0.10 U							
2-Methylphenol	95-48-7	800							10 U		
2-Nitrophenol	88-75-5								10 U		
3- & 4-Methylphenol	15831-10-4	800							10 U		
4,6-Dinitro-o-cresol	534-52-1	1.3							30 U		
4-Chloro-3-methylphenol	59-50-7	1,600							Į		
4-Nitrophenol	100-02-7	1,600							10 U		
Acenaphthene	83-32-9	480		0.10 U					30		
Acenaphthylene	208-96-8			0.10 U							
Acrylonitrile	107-13-1				10 U	10 U	10 U			10 U	
Anthracene	120-12-7	2,400		0.10 U					1.3		
Benzo(a)anthracene	56-55-3		0.10 U	0.10 U					0.10 U		
Benzo(a)pyrene	50-32-8		0.10 U	0.10 U					0.10 U		
Benzo(b)fluoranthene	205-99-2		0.10 U	0.10 U					0.10 U		
Benzo(g,h,i)perylene	191-24-2			0.10 U							
Benzo(k)fluoranthene	207-08-9		0.10 U	0.10 U					0.10 U		
Benzoic acid	65-85-0								100 U		
Benzyl alcohol	100-51-6								1.0 U		
Bis(2-chloroethoxy)methane	111-91-1								1.0 U		
Bis(2-ethylhexyl)phthalate	117-81-7										
Butyl benzyl phthalate	85-68-7								1.0 U		
Carbazole	86-74-8								2.8		
Chrysene	218-01-9		0.10 U	0.10 U					0.10 U		
Dibenzo(a,h)anthracene	53-70-3		0.10 U	0.10 U					0.10 U		
Indeno(1,2,3-c,d)pyrene	193-39-5		0.10 U	0.10 U					0.10 U		
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	0.10 U	0.10 U					0.10 U		
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	0.10 U	0.10 U					0.10 U		
Dibenzofuran	132-64-9								13		
Diethylphthalate	84-66-2					İ			1.0 U	1	
Dimethyl phthalate	131-11-3								1		
Di-n-butyl phthalate	84-74-2								1.0 U		
Di-n-octyl phthalate	117-84-0								1.0 0		
Fluoranthene	206-44-0	640		0.10 U					1.8		
Fluorene	86-73-7	320		0.10 U		 		1	6.5	 	
Hexachlorobutadiene	87-68-3			1.0 U	2.0 U	2.0 U	2.0 U	1	1.0 U	2.0 U	
Hexachlorocyclopentadiene	77-47-4			1.00	2.00	2.00	2.00		1.00	2.00	
Isophorone	78-59-1					 		1	1.0 U	 	
Naphthalene	91-20-3		0 77 (3)	(4)	(3)	(3)	(3)	}	5.1	(3)	
		160 (3)	0.77 (3)	0.10 U ⁽⁴⁾	2.0 U ⁽³⁾	2.0 U ⁽³⁾	2.0 U ⁽³⁾	<u> </u>		4.8 (3)	
N-Nitroso-di-n-propylamine	621-64-7								1.0 U		
N-Nitrosodiphenylamine	86-30-6	-							1.0 U		
Pentachlorophenol	87-86-5	1.0				1			10 U	1	



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name			MW-01	veli Analyticai Data 2	MW-02	l	MW-04		
		Sample Name	WCD-MW1-38	WCD-MW01-112712	WCD-MW01-030513	WCD-MW01-060413	WCD-MW01-100113	MW02-WL-03292021	WCD-MW4-33	MW04-WL-03292021	
		Sample Date	6/22/2010	11/27/2012	3/5/2013	6/4/2013	10/1/2013	3/29/2021	6/21/2010	WCD-MW04-120511 12/5/2011	3/29/2021
Analyte	CAS No.	Screening Level (1)	2,22,222		5/1/2020	5, ,,====	,-,	-,,	0,00,000		2,20,202
Semivolatile Organic Compounds (cor		Screening Level									
Phenanthrene	85-01-8			0.10 U					4.6		
Phenol	108-95-2	4,800							10 U		
Pyrene	129-00-0	240		0.10 U					1.2		
Total Naphthalenes		160		0.10 U							
Volatile Organic Compounds											
1,1,1,2-Tetrachloroethane	630-20-6			1.0 U	2.0 U	2.0 U	2.0 U		l e	2.0 U	
1,1,1-Trichloroethane	71-55-6	200		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
1,1,2,2-Tetrachloroethane	79-34-5			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
1,1,2-Trichloroethane	79-00-5			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
1,1-Dichloroethane	75-34-3	7.7		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
1,1-Dichloroethene	75-35-4			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
1,1-Dichloropropene	563-58-6			1.0 U	2.0 U	2.0 U	2.0 U		1	2.0 U	
1,2,3-Trichlorobenzene	87-61-6			1.0 U	2.0 U	2.0 U	2.0 U		1	2.0 U	
1,2,3-Trichloropropane	96-18-4			1.0 U	2.0 U	2.0 U	2.0 U		1	2.0 U	
1,2,4-Trichlorobenzene	120-82-1			1.0 U	2.0 U	2.0 U	2.0 U		1.0 U	2.0 U	
1,2,4-Trimethylbenzene	95-63-6	80		1.0 U	2.0 U	2.0 U	2.0 U		1.0 0	2.0 U	
1,2-Dibromo-3-chloropropane	96-12-8			1.0 U	10 U	10 U	10 U		1	10 U	
1,2-Dibromoethane	106-93-4	-		1.0 U	0.010 U	0.010 U	0.010 U			0.010 U	
1,2-Dichlorobenzene	95-50-1			1.0 U	2.0 U	2.0 U	2.0 U		1.0 U	2.0 U	
1,2-Dichloroethane	107-06-2			1.0 U	2.0 U	2.0 U	2.0 U		1.0 0	2.0 U	
1,2-Dichloropropane	78-87-5			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
1,3,5-Trimethylbenzene	108-67-8	80		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
1,3-Dichlorobenzene	541-73-1			1.0 U	2.0 U	2.0 U	2.0 U		1.0 U	2.0 U	
1,3-Dichloropropane	142-28-9			1.0 U	2.0 U	2.0 U	2.0 U		1.0 0	2.0 U	
1,4-Dichlorobenzene	106-46-7			1.0 U	2.0 U	2.0 U	2.0 U		1.0 U	2.0 U	
2,2-Dichloropropane	594-20-7			1.0 U	2.0 U	2.0 U	2.0 U		1.0 0	2.0 U	
2,4-Dinitrotoluene	121-14-2			1.00	2.0 0	2.00	2.00		1.0 U	2.00	
2,6-Dinitrotoluene	606-20-2								1.0 0		
2-Chlorotoluene	95-49-8			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
2-Hexanone	591-78-6			1.0 U	10 U	10 U	10 U			10 U	
2-Nitroaniline	88-74-4			1.00	10 0	10 0	10 0		1	10 0	
3-Nitroaniline	99-09-2								3.0 U		
4-Chloroaniline	106-47-8								3.0 0		
4-Chlorophenyl phenyl ether	7005-72-3								1.0 U		
4-Chlorotoluene	106-43-4			1.0 U	2.0 U	2.0 U	2.0 U		1.00	2.0 U	
4-Nitroaniline	100-01-6			1.00	2.0 0	2.00	2.00		10 U	2.00	
Acetone	67-64-1			10 U	25 U	25 U	25 U		10 0	25 U	
Benzene	71-43-2	5.0		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Bis(2-chloroethyl)ether	111-44-4	3.0		1.00	2.00	2.00	2.00		1.0 U	2.00	
Bis(2-chloroisopropyl)ether	39638-32-9			<u> </u>		<u> </u>		1	1.0 U	 	
Bromobenzene	108-86-1			1.0 U	2.0 U	2.0 U	2.0 U	1	1.00	2.0 U	
Bromochloromethane	74-97-5			1.0 U	2.0 U	2.0 U	2.0 U	}	1	2.0 U	
Bromodichloromethane	75-27-4			1.0 U	2.0 U	2.0 U	2.0 U		1	2.0 U	
Bromoform	75-25-2			1.0 U	2.0 U	2.0 U	2.0 U	1	1	2.0 U	
Bromomethane	75-25-2			1.0 U	2.0 U	2.0 U	2.0 U	}	1	2.0 U	
Carbon disulfide	75-15-0			1.00	2.0 U	2.0 U	2.0 U	}	1	2.0 U	
Carbon disulfide Carbon tetrachloride	75-13-0 56-23-5			1.0 U	2.0 U	2.0 U	2.0 U		1	2.0 U	
Chlorobenzene	108-90-7			1.0 U	2.0 U	2.0 U	2.0 U	1	1	2.0 U	
Chloroethane	75-00-3			1.0 U	2.0 U	2.0 U	2.0 U	}	1	2.0 U	
Chloroform	67-66-3	70		1.0 U	2.0 U	2.0 U	2.0 U		1	2.0 U	



Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (µg/L)

Loc	ation Name				MW-01			MW-02		MW-04	
		Sample Name	WCD-MW1-38	WCD-MW01-112712	WCD-MW01-030513	WCD-MW01-060413	WCD-MW01-100113	MW02-WL-03292021	WCD-MW4-33	WCD-MW04-120511	MW04-WL-03292021
		Sample Date	6/22/2010	11/27/2012	3/5/2013	6/4/2013	10/1/2013	3/29/2021	6/21/2010	12/5/2011	3/29/2021
Analyte	CAS No.	Screening Level (1)									
Volatile Organic Compounds (cont.)	•			•					-	•	
Chloromethane	74-87-3			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
cis-1,2-Dichloroethene	156-59-2	70		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
cis-1,3-Dichloropropene	10061-01-5			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Cymene	99-87-6			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Dibromochloromethane	124-48-1	-		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Dibromomethane	74-95-3			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Dichlorodifluoromethane	75-71-8	-		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Ethylbenzene	100-41-4	700		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Hexachlorobenzene	118-74-1								1.0 U		
Hexachloroethane	67-72-1								1.0 U		
Isopropylbenzene	98-82-8	800		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Methyl ethyl ketone	78-93-3			10 U	10 U	10 U	10 U			10 U	
Methyl isobutyl ketone	108-10-1	640		1.0 U	10 U	10 U	10 U			10 U	
Methylene chloride	75-09-2			1.0 U	5.0 U	5.0 U	5.0 U			5.0 U	
Methyl-tert-butyl ether	1634-04-4			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
n-Butylbenzene	104-51-8			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Nitrobenzene	98-95-3								1.0 U		
n-Propylbenzene	103-65-1	800		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
sec-Butylbenzene	135-98-8	-		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Styrene	100-42-5	100	1.0 U	1.0 U	2.0 U	2.0 U	2.0 U		1.0 U	2.0 U	
tert-Butylbenzene	98-06-6			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Tetrachloroethene	127-18-4			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Toluene	108-88-3	1,000		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
trans-1,2-Dichloroethene	156-60-5			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
trans-1,3-Dichloropropene	10061-02-6			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Trichloroethene	79-01-6	5.0		1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Trichlorofluoromethane	75-69-4			1.0 U	2.0 U	2.0 U	2.0 U			2.0 U	
Vinyl chloride	75-01-4			0.20 U	0.20 U	0.20 U	0.20 U			0.20 U	
Xylene (meta & para)	108-38-3/106-42-3				4.0 U	4.0 U	4.0 U			4.0 U	
Xylene (ortho)	95-47-6				2.0 U	2.0 U	2.0 U			2.0 U	
Xylene (total)	1330-20-7	1,000		1.0 U	4.0 U	4.0 U	4.0 U			4.0 U	
Other											
PBDE-003	101-55-3								1.0 U		

Blank cells are intentional

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

- Not available or not established.

The screening level is a proposed Site cleanup level.

- 1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from the February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.
- 2 A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest RL was selected. 3 Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level for total naphthalenes. 3 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level. 5 Xylene (total) as calculated by Floyd | Snider was preferred over laboratory reported result.

Abbreviations:

CAS = Chemical Abstracts Service

PBDE = Polybrominated diphenyl ethers TEQ = Toxic equivalent

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

USEPA = U.S. Environmental Protection Agency

μg/L = Micrograms per liter MTCA = Model Toxics Control Act

Qualifiers:

- J Analyte was detected and the concentration is estimated.
- U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

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		Location Name	MW-05									
		Sample Name	WCD-MW5-32	WCD-MW05-120511	WCD-MW05-112712	WCD-MW05-030513	WCD-MW05-060413	WCD-MW05-100113	MW05-09152020	MW05-WL-03292021		
		Sample Date	6/21/2010	12/5/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013	9/15/2020	3/29/2021		
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds		ourcening acres										
1-Methylnaphthalene	90-12-0				0.10 U				0.10 U			
2,4,5-Trichlorophenol	95-95-4	1,600			3.23				2.0 U			
2,4,6-Trichlorophenol	88-06-2	8.0							2.0 U			
2.4-Dichlorophenol	120-83-2	48							2.0 U			
2,4-Dimethylphenol	105-67-9	320							1.0 U			
2,4-Dinitrophenol	51-28-5	32							2.0 UJ			
2-Chloronaphthalene	91-58-7								, ,,			
2-Chlorophenol	95-57-8	40							1.0 U			
2-Methylnaphthalene	91-57-6				0.80				0.10 U			
2-Methylphenol	95-48-7	800			0.00				1.0 U			
2-Nitrophenol	88-75-5								2.0 U			
3- & 4-Methylphenol	15831-10-4	800				<u> </u>		<u> </u>	1.0 U			
4,6-Dinitro-o-cresol	534-52-1	1.3							5.0 UJ			
4-Chloro-3-methylphenol	59-50-7	1,600							5.0 U			
4-Nitrophenol	100-02-7	1,600						+	5.0 UJ			
Acenaphthene	83-32-9	480			0.90			+	0.20 U			
Acenaphthylene	208-96-8	460			0.10 U			1	0.20 U			
				10.11	0.10 0	10.11	10.11	10.11	0.20 0			
Acrylonitrile	107-13-1 120-12-7	2,400		10 U	0.10 U	10 U	10 U	10 U	0.10 U			
Anthracene	56-55-3		2.7		0.10 U				0.10 U			
Benzo(a)anthracene		-	2.7 1.6		0.10 U							
Benzo(a)pyrene	50-32-8		2.4		0.10 U				0.10 U			
Benzo(b)fluoranthene	205-99-2		2.4						0.20 U			
Benzo(g,h,i)perylene	191-24-2		0.59		0.10 U				0.20 U			
Benzo(k)fluoranthene	207-08-9		0.59		0.10 U				0.20 U			
Benzoic acid	65-85-0											
Benzyl alcohol	100-51-6											
Bis(2-chloroethoxy)methane	111-91-1											
Bis(2-ethylhexyl)phthalate	117-81-7											
Butyl benzyl phthalate	85-68-7											
Carbazole	86-74-8											
Chrysene	218-01-9		3.5		0.10 U				0.20 U			
Dibenzo(a,h)anthracene	53-70-3		0.17		0.10 U				0.20 U			
Indeno(1,2,3-c,d)pyrene	193-39-5		0.62		0.10 U				0.10 U			
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	2.3		0.10 U				0.10 U			
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	2.3		0.10 U				0.10 U			
Dibenzofuran	132-64-9											
Diethylphthalate	84-66-2											
Dimethyl phthalate	131-11-3											
Di-n-butyl phthalate	84-74-2											
Di-n-octyl phthalate	117-84-0											
Fluoranthene	206-44-0	640			0.10 U				0.10 U			
Fluorene	86-73-7	320			0.10 U				0.10 U			
Hexachlorobutadiene	87-68-3		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Hexachlorocyclopentadiene	77-47-4											
Isophorone	78-59-1											
Naphthalene	91-20-3	160 ⁽³⁾	0.39 ⁽³⁾	31 ⁽³⁾	11 (4)	70 ⁽³⁾	5.0 ⁽³⁾	2.0 U ⁽³⁾	2.0 U ⁽³⁾			
N-Nitroso-di-n-propylamine	621-64-7											
N-Nitrosodiphenylamine	86-30-6											
Pentachlorophenol	87-86-5	1.0							0.10 UJ			



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name					MW-05			
		Sample Name	WCD-MW5-32	WCD-MW05-120511	WCD-MW05-112712	WCD-MW05-030513	WCD-MW05-060413	WCD-MW05-100113	MW05-09152020	MW05-WL-03292021
		Sample Date	6/21/2010	12/5/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013	9/15/2020	3/29/2021
Analyte	CAS No.	Screening Level (1)		,		.,.,	.,,			
Semivolatile Organic Compounds (co		Screening Level								
Phenanthrene	85-01-8				0.10 U				0.10 U	
Phenol	108-95-2	4,800							2.0 UJ	
Pyrene	129-00-0	240			0.10 U				0.20 U	
Total Naphthalenes	-	160			12				2.0 U	
Volatile Organic Compounds										
1,1,1,2-Tetrachloroethane	630-20-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,1,1-Trichloroethane	71-55-6	200	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	
1,1,2,2-Tetrachloroethane	79-34-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,1,2-Trichloroethane	79-00-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,1-Dichloroethane	75-34-3	7.7	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	
1,1-Dichloroethene	75-35-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,1-Dichloropropene	563-58-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,2,3-Trichlorobenzene	87-61-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,2,3-Trichloropropane	96-18-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,2,4-Trichlorobenzene	120-82-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,2,4-Trimethylbenzene	95-63-6	80	1.1	2.0 U	1.0 U	2.5	2.0 U	2.0 U	1.0 U	
1,2-Dibromo-3-chloropropane	96-12-8		10 U	10 U	1.0 U	10 U	10 U	10 U		
1,2-Dibromoethane	106-93-4		1.0 U	0.010 U	1.0 U	0.010 U	0.010 U	0.010 U		
1,2-Dichlorobenzene	95-50-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,2-Dichloroethane	107-06-2		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,2-Dichloropropane	78-87-5	_	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,3,5-Trimethylbenzene	108-67-8	80	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	
1,3-Dichlorobenzene	541-73-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,3-Dichloropropane	142-28-9	_	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
1,4-Dichlorobenzene	106-46-7		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
2,2-Dichloropropane	594-20-7		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
2,4-Dinitrotoluene	121-14-2									
2,6-Dinitrotoluene	606-20-2									
2-Chlorotoluene	95-49-8		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
2-Hexanone	591-78-6		10 U	10 U	1.0 U	10 U	10 U	10 U		
2-Nitroaniline	88-74-4									
3-Nitroaniline	99-09-2									
4-Chloroaniline	106-47-8									
4-Chlorophenyl phenyl ether	7005-72-3									
4-Chlorotoluene	106-43-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
4-Nitroaniline	100-01-6									
Acetone	67-64-1		10 U	25 U	10 U	25 U	25 U	25 U		
Benzene	71-43-2	5.0	0.35 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	
Bis(2-chloroethyl)ether	111-44-4									
Bis(2-chloroisopropyl)ether	39638-32-9									
Bromobenzene	108-86-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Bromochloromethane	74-97-5			2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Bromodichloromethane	75-27-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Bromoform	75-25-2		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Bromomethane	74-83-9	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Carbon disulfide	75-15-0			2.0 U		2.0 U	2.0 U	2.0 U		
Carbon tetrachloride	56-23-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Chlorobenzene	108-90-7	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Chloroethane	75-00-3		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		
Chloroform	67-66-3	70	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	



Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (µg/L)

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		Location Name	MW-05									
		Sample Name	WCD-MW5-32	WCD-MW05-120511	WCD-MW05-112712	WCD-MW05-030513	WCD-MW05-060413	WCD-MW05-100113	MW05-09152020	MW05-WL-03292021		
		Sample Date	6/21/2010	12/5/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013	9/15/2020	3/29/2021		
Analyte	CAS No.	Screening Level (1)										
Volatile Organic Compounds (cont.)												
Chloromethane	74-87-3	-	10 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	2.0 U			
cis-1,2-Dichloroethene	156-59-2	70	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U			
cis-1,3-Dichloropropene	10061-01-5	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Cymene	99-87-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Dibromochloromethane	124-48-1	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Dibromomethane	74-95-3		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Dichlorodifluoromethane	75-71-8	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Ethylbenzene	100-41-4	700	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U			
Hexachlorobenzene	118-74-1											
Hexachloroethane	67-72-1											
Isopropylbenzene	98-82-8	800	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U			
Methyl ethyl ketone	78-93-3		10 U	10 U	10 U	10 U	10 U	10 U				
Methyl isobutyl ketone	108-10-1	640	10 U	10 U	1.0 U	10 U	10 U	10 U	1.3 U			
Methylene chloride	75-09-2		5.0 U	5.0 U	1.0 U	5.0 U	5.0 U	5.0 U				
Methyl-tert-butyl ether	1634-04-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
n-Butylbenzene	104-51-8			2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Nitrobenzene	98-95-3	-										
n-Propylbenzene	103-65-1	800	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U			
sec-Butylbenzene	135-98-8		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Styrene	100-42-5	100	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U			
tert-Butylbenzene	98-06-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Tetrachloroethene	127-18-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Toluene	108-88-3	1,000	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U			
trans-1,2-Dichloroethene	156-60-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
trans-1,3-Dichloropropene	10061-02-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Trichloroethene	79-01-6	5.0	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	0.50 U			
Trichlorofluoromethane	75-69-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Vinyl chloride	75-01-4		0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U				
Xylene (meta & para)	108-38-3/106-42-3		2.0 U	4.0 U		4.0 U	4.0 U	4.0 U	1.0 U			
Xylene (ortho)	95-47-6		1.0 U	2.0 U		2.0 U	2.0 U	2.0 U	1.0 U			
Xylene (total)	1330-20-7	1,000	2.0 U	4.0 U	1.0 U	4.0 U	4.0 U	4.0 U	1.0 U ⁽⁵⁾			
Other												
PBDE-003	101-55-3											

Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

-- Not available or not established.

BOLD Detected result exceeds criteria. The screening level is a proposed Site cleanup level.

- 1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.

 2 A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest Rt. was selected. 3 Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level 5. Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level. 5 Windows analyzed by the March 10 of 10 o

Abbreviations:

CAS = Chemical Abstracts Service

PBDE = Polybrominated diphenyl ethers TEQ = Toxic equivalent

cPAH = Carcinogenic polycyclic aromatic hydrocarbon μg/L = Micrograms per liter

USEPA = U.S. Environmental Protection Agency

MTCA = Model Toxics Control Act

Qualifiers:

J Analyte was detected and the concentration is estimated.

U Analyte is not detected at the associated reporting limit.

Modified from Floyd | Snider 2022 draft RI/FS.

UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Modified from Floyd | Snider 2022 draft RI/FS.

Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name					MW-06	5				
		Sample Name	WCD-MW6-31	WCD-MW06-120611	WCD-MW06-112712	WCD-MW06-030513	WCD-MW06-060413	WCD-MW06-100113	MW06-09152020 ⁽²⁾	MW-06-12172020	MW06-032921	MW06-06222021
		Sample Date	6/21/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013	9/15/2020	12/17/2020	3/29/2021	6/22/2021
Analyte	CAS No.	Screening Level (1)	0,11,111	, _, _,	,,	5/5/2020	5, 4,2020	,	0,10,100		5,25,2522	1,12,202
Semivolatile Organic Compounds		Screening Level										
1-Methylnaphthalene	90-12-0				800		470	480	570	400	370	550
2,4,5-Trichlorophenol	95-95-4	1,600	10 U		000			100	2.0 U	2.0 UJ	3,0	330
2,4,6-Trichlorophenol	88-06-2	8.0	10 U						2.0 U	2.0 UJ		
2,4-Dichlorophenol	120-83-2	48	10 U						2.0 U	2.0 UJ		
2,4-Dimethylphenol	105-67-9	320	10 U						0.98 U	0.99 UJ		
2,4-Dinitrophenol	51-28-5	32	30 U						2.0 UJ	2.0 UJ		
2-Chloronaphthalene	91-58-7		1.0 U						2.0 03	2.0 03		
2-Chlorophenol	95-57-8	40	10 U						0.98 U	0.99 UJ		
·	91-57-6		780		1,000		710	980	740	550	460	720
2-Methylnaphthalene	95-48-7	800	780 10 U		1,000		710	960	1.5	330	460	720
2-Methylphenol	88-75-5	800	10 U						2.0 U	2.0 UJ		
2-Nitrophenol	15831-10-4		10 U						2.0 0	0.99 UJ		
3- & 4-Methylphenol		800										
4,6-Dinitro-o-cresol	534-52-1	1.3	30 U						49 U	5.0 UJ	1	
4-Chloro-3-methylphenol	59-50-7	1,600	10 U						4.9 U	5.0 UJ		
4-Nitrophenol	100-02-7	1,600	10 U						4.9 UJ	5.0 UJ		
Acenaphthene	83-32-9	480	220		6.9		170	190	190	110	140	160
Acenaphthylene	208-96-8	-	20		240		15	22	19	9.9 J		11
Acrylonitrile	107-13-1			10 U		10 U	10 U	10 U				
Anthracene	120-12-7	2,400	19		7.6		6.0	20 U	8.4	4.3	6.0	4.6
Benzo(a)anthracene	56-55-3		6.9		0.10 U	0.75	0.74	0.68	1.8	0.57	1.6	1.1
Benzo(a)pyrene	50-32-8		5.9		0.10 U	0.48	0.52	0.49	1.3	0.16	1.4	0.67
Benzo(b)fluoranthene	205-99-2		7.8		0.10 U	0.45	0.55	0.45	1.1	0.26	1.2	0.60
Benzo(g,h,i)perylene	191-24-2		2.2		0.10 U		0.23	0.61	0.47	0.50 U		0.37
Benzo(k)fluoranthene	207-08-9		5.0 U		0.10 U	0.35	0.54	0.31	0.85	0.22	1.1	0.56
Benzoic acid	65-85-0		100 U									
Benzyl alcohol	100-51-6		1.0 U									
Bis(2-chloroethoxy)methane	111-91-1		1.0 U									
Bis(2-ethylhexyl)phthalate	117-81-7		10 U									
Butyl benzyl phthalate	85-68-7		1.0 U									
Carbazole	86-74-8	-	110									
Chrysene	218-01-9		6.8		0.90	0.77	0.78	0.69	1.3	0.40	1.2	0.80
Dibenzo(a,h)anthracene	53-70-3		5.0 U		0.10 U	0.29	0.080 U	0.44	0.22	0.14	0.40 U	0.40 U
Indeno(1,2,3-c,d)pyrene	193-39-5		5.0 U		0.10 U	0.35	0.16	0.54	0.44	0.16	0.58	0.33
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	8.2		0.084	0.71	0.73	0.74	1.8	0.30	1.9	0.96
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	7.4		0.0090	0.71	0.73	0.74	1.8	0.30	1.9	0.94
Dibenzofuran	132-64-9		69									
Diethylphthalate	84-66-2		1.0 U									
Dimethyl phthalate	131-11-3		1.0 U						1			
Di-n-butyl phthalate	84-74-2		1.0 U									
Di-n-octyl phthalate	117-84-0		1.0 U									
Fluoranthene	206-44-0	640	31		11		6.3	7.0	10	5.8		8.2
Fluorene	86-73-7	320	89		54		35	49	60	50 U		54
Hexachlorobutadiene	87-68-3		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U		30 0		J-
Hexachlorocyclopentadiene	77-47-4		3.0 U	2.00	1.00	2.00	2.00	2.00	 			
	78-59-1		1.0 U						-			
Isophorone Naphthalene	91-20-3		5,900	(3)	(4)	(3)	(4)	(4)	5,700	2,900	3,700	(4)
		160 (3)	·	9,200 (3)	9,800 ⁽⁴⁾	1,100 (3)	7,400 ⁽⁴⁾	9,900 ⁽⁴⁾	3,700	2,900	3,700	5,200 ⁽⁴⁾
N-Nitroso-di-n-propylamine	621-64-7		1.0 U						 			
N-Nitrosodiphenylamine	86-30-6		1.0 U									
Pentachlorophenol	87-86-5	1.0	10 U	1	1	1	I	1	0.098 UJ	0.20 UJ	0.27 U	0.20 U



Modified from Floyd | Snider 2022 draft RI/FS.

Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name					MW-06	i				
		Sample Name	WCD-MW6-31	WCD-MW06-120611	WCD-MW06-112712	WCD-MW06-030513	WCD-MW06-060413	WCD-MW06-100113	MW06-09152020 ⁽²⁾	MW-06-12172020	MW06-032921	MW06-06222021
		Sample Date	6/21/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013	9/15/2020	12/17/2020	3/29/2021	6/22/2021
Analyte	CAS No.	Screening Level (1)	0,21,2010	12/0/2011	12/2//2012	3,3,2013	0,4,2020	10/1/2010	3/13/2020	11/1// 1010	5,25,2022	0/12/2021
Semivolatile Organic Compounds (con		Screening Level		1							l	
Phenanthrene	85-01-8		110		26		22	22	42	50 U	I	35
Phenol	108-95-2	4,800	10 U		20				2.0 UJ	2.0 UJ		33
Pyrene	129-00-0	240	30		7.1		4.9	4.2	7.5	3.8		5.2
Total Naphthalenes		160	6,700		12,000		8,600	11,000	7,000	3,800	4,500	6,500
			-,,			Organic Compounds	5,222	,	-,	-,	.,	-,
1,1,1,2-Tetrachloroethane	630-20-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U			I	
1.1.1-Trichloroethane	71-55-6	200	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	4.0 U	4.0 U
1,1,2,2-Tetrachloroethane	79-34-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,1,2-Trichloroethane	79-00-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,1-Dichloroethane	75-34-3	7.7	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	5.0 U	5.0 U
1,1-Dichloroethene	75-35-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				0.00
1,1-Dichloropropene	563-58-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,2,3-Trichlorobenzene	87-61-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U			1	
1,2,3-Trichloropropane	96-18-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,2,4-Trichlorobenzene	120-82-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,2,4-Trimethylbenzene	95-63-6	80	91 J	150	170	160	150	170	160	120	97	110
1,2-Dibromo-3-chloropropane	96-12-8		10 U	10 U	1.0 U	10 U	10 U	10 U				
1,2-Dibromoethane	106-93-4		1.0 U	0.010 U	1.0 U	0.010 U	0.010 U	0.010 U				
1,2-Dichlorobenzene	95-50-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,2-Dichloroethane	107-06-2		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,2-Dichloropropane	78-87-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,3,5-Trimethylbenzene	108-67-8	80	51	51	63	57	54	65	47	41	31	36
1,3-Dichlorobenzene	541-73-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	.,		0.1	30
1,3-Dichloropropane	142-28-9		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
1,4-Dichlorobenzene	106-46-7		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
2,2-Dichloropropane	594-20-7		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
2,4-Dinitrotoluene	121-14-2		1.0 U	2.00	1.00	2.00	2.00	2.00				
2,6-Dinitrotoluene	606-20-2		1.0 U									
2-Chlorotoluene	95-49-8		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
2-Hexanone	591-78-6		10 U	10 U	1.0 U	10 U	10 U	10 U				
2-Nitroaniline	88-74-4		1.0 U	10 0	1.00	10 0	10 0	10 0				
3-Nitroaniline	99-09-2		3.0 U									
4-Chloroaniline	106-47-8		3.0 U									
4-Chlorophenyl phenyl ether	7005-72-3		1.0 U									
4-Chlorotoluene	106-43-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
4-Nitroaniline	100-01-6		10 U									
Acetone	67-64-1		10 U	25 U	10 U	25 U	25 U	25 U				
Benzene	71-43-2	5.0	0.35 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	4.4 U	4.4 U
Bis(2-chloroethyl)ether	111-44-4		1.0 U			=:= =						
Bis(2-chloroisopropyl)ether	39638-32-9		1.0 U									
Bromobenzene	108-86-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Bromochloromethane	74-97-5			2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Bromodichloromethane	75-27-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Bromoform	75-25-2		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Bromomethane	74-83-9		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Carbon disulfide	75-15-0			2.0 U		2.0 U	2.0 U	2.0 U				
Carbon tetrachloride	56-23-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Chlorobenzene	108-90-7	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Chloroethane	75-00-3		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Chloroform	67-66-3	70	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	5.0 U	5.0 U
C.11010101111	0, 00 3	,,	1.0 0	2.0 0	1.0 0	2.0 0	2.0 0	2.0 0	1.0 0	1.0 0	5.0 0	3.0 0



Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (μg/L)

		Location Name					MW-06	i				
		Sample Name	WCD-MW6-31	WCD-MW06-120611	WCD-MW06-112712	WCD-MW06-030513	WCD-MW06-060413	WCD-MW06-100113	MW06-09152020 ⁽²⁾	MW-06-12172020	MW06-032921	MW06-06222021
		Sample Date	6/21/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013	9/15/2020	12/17/2020	3/29/2021	6/22/2021
Analyte	CAS No.	Screening Level (1)										
Volatile Organic Compounds (cont.)												
Chloromethane	74-87-3		10 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	7.5 U	7.5 U
cis-1,2-Dichloroethene	156-59-2	70	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	5.0 U	5.0 U
cis-1,3-Dichloropropene	10061-01-5	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Cymene	99-87-6	-	6.2	6.4	5.2	6.3	5.1	6.1				
Dibromochloromethane	124-48-1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Dibromomethane	74-95-3	-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Dichlorodifluoromethane	75-71-8		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Ethylbenzene	100-41-4	700	7.6	2.1	3.0	2.1	2.0 U	3.1	2.4	1.5	4.0 U	4.0 U
Hexachlorobenzene	118-74-1		1.0 U									
Hexachloroethane	67-72-1	-	1.0 U									
Isopropylbenzene	98-82-8	800	7.5	6.0	7.1	7.1	5.5	6.5	7.6	5.6	5.0 U	5.7
Methyl ethyl ketone	78-93-3		10 U	10 U	10 U	10 U	10 U	10 U			15 U	
Methyl isobutyl ketone	108-10-1	640	10 U	10 U	1.0 U	10 U	10 U	10 U	1.3 U	1.3 U		13 U
Methylene chloride	75-09-2		5.0 U	5.0 U	1.0 U	5.0 U	5.0 U	5.0 U				
Methyl-tert-butyl ether	1634-04-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
n-Butylbenzene	104-51-8			2.0 U	8.6	2.0 U	2.0 U	2.0 U				
Nitrobenzene	98-95-3		1.0 U									
n-Propylbenzene	103-65-1	800	4.7	3.2	5.1	2.0 U	2.0 U	4.4	6.3	4.6	5.0 U	5.0 U
sec-Butylbenzene	135-98-8		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Styrene	100-42-5	100	9.2	3.5	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	5.0 U	5.0 U
tert-Butylbenzene	98-06-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Tetrachloroethene	127-18-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Toluene	108-88-3	1,000	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	7.5 U	7.5 U
trans-1,2-Dichloroethene	156-60-5		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
trans-1,3-Dichloropropene	10061-02-6		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Trichloroethene	79-01-6	5.0	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U	0.50 U	0.50 U	5.0 U	5.0 U
Trichlorofluoromethane	75-69-4		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U				
Vinyl chloride	75-01-4		0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U				
Xylene (meta & para)	108-38-3/106-42-3		56	27		23	16	21	17	11	10 U	14
Xylene (ortho)	95-47-6		48	25		19	14	18	11	6.9	5.0 U	8.1
Xylene (total)	1330-20-7	1,000	100	52	63	42	30	39	28 ⁽⁵⁾	18	10 U ⁽⁵⁾	22 ⁽⁵⁾
Other											-	-
PBDE-003	101-55-3		1.0 U									

Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

-- Not available or not established.

BOLD Detected result exceeds criteria.

The screening level is a proposed Site cleanup level.

- 1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.

 2 A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest Rt. was selected. 3 Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level 5. Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level. 5 Windows analyzed by the March 10 of 10 o

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

μg/L = Micrograms per liter MTCA = Model Toxics Control Act PBDE = Polybrominated diphenyl ethers

TEQ = Toxic equivalent

USEPA = U.S. Environmental Protection Agency

Qualifiers:

J Analyte was detected and the concentration is estimated.

Modified from Floyd | Snider 2022 draft RI/FS.

- U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name			MW-07	, , , , , , , , , , , , , , , , , , , ,		MW-08			MW-09	
			WCD-MW7-36	WCD-MW07-120511	WCD-MW07-112712	WCD-MW07-100113	MW07-WL-03292021	WCD-MW8-37	WCD-MW9-65	WCD-MW09-120511	WCD-MW09-112712	WCD-MW09-100113
		Sample Name Sample Date	6/21/2010	12/5/2011	11/27/2012	10/1/2013	3/29/2021	6/22/2010		12/5/2011	11/27/2012	10/1/2013
A b. d	CACAL		6/21/2010	12/5/2011	11/2//2012	10/1/2015	3/23/2021	6/22/2010	6/21/2010	12/3/2011	11/2//2012	10/1/2013
Analyte	CAS No.	Screening Level (1)						<u> </u>				1
Semivolatile Organic Compounds 1-Methylnaphthalene	90-12-0			l	0.10 U	l	l	ı			0.10 U	0.023
2,4,5-Trichlorophenol	95-95-4	1,600			0.10 0			1			0.10 0	0.023
2,4,6-Trichlorophenol	88-06-2	8.0						1				
2,4-Dichlorophenol	120-83-2	48						-				
2,4-Dimethylphenol	105-67-9	320						1				
2,4-Dinitrophenol	51-28-5	320						-				
	91-58-7							-				
2-Chloronaphthalene 2-Chlorophenol	95-57-8	40										
	95-57-8	40			1.6			-			0.1011	0.020.11
2-Methylnaphthalene					1.6						0.10 U	0.020 U
2-Methylphenol	95-48-7	800										
2-Nitrophenol	88-75-5											+
3- & 4-Methylphenol	15831-10-4	800						1			1	
4,6-Dinitro-o-cresol	534-52-1	1.3						1			1	
4-Chloro-3-methylphenol	59-50-7	1,600										
4-Nitrophenol	100-02-7	1,600										
Acenaphthene	83-32-9	480			1.2						0.10 U	0.12
Acenaphthylene	208-96-8				2.4						0.10 U	0.020 U
Acrylonitrile	107-13-1			10 U		10 U				10 U		10 U
Anthracene	120-12-7	2,400			0.10 U						0.10 U	0.039
Benzo(a)anthracene	56-55-3		0.96		0.10 U			0.10 U	0.14		0.10 U	0.021
Benzo(a)pyrene	50-32-8		0.43		0.10 U			0.10 U	0.10 U		0.10 U	0.029 U
Benzo(b)fluoranthene	205-99-2		0.58		0.10 U			0.10 U	0.10 U		0.10 U	0.020 U
Benzo(g,h,i)perylene	191-24-2				0.10 U						0.10 U	0.077
Benzo(k)fluoranthene	207-08-9		0.21		0.10 U			0.10 U	0.10 U		0.10 U	0.020 U
Benzoic acid	65-85-0											
Benzyl alcohol	100-51-6											
Bis(2-chloroethoxy)methane	111-91-1											
Bis(2-ethylhexyl)phthalate	117-81-7	-										
Butyl benzyl phthalate	85-68-7	-										
Carbazole	86-74-8											
Chrysene	218-01-9		0.93		0.10 U			0.10 U	0.11		0.10 U	0.020 U
Dibenzo(a,h)anthracene	53-70-3		0.10 U		0.10 U			0.10 U	0.10 U		0.10 U	0.070
Indeno(1,2,3-c,d)pyrene	193-39-5		0.18		0.10 U			0.10 U	0.10 U		0.10 U	0.072
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	0.64		0.10 U			0.10 U	0.085		0.10 U	0.033
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	0.63		0.10 U			0.10 U	0.015		0.10 U	0.016
Dibenzofuran	132-64-9							1				
Diethylphthalate	84-66-2							1			†	
Dimethyl phthalate	131-11-3							1				
Di-n-butyl phthalate	84-74-2							1				
Di-n-octyl phthalate	117-84-0							 				
Fluoranthene	206-44-0	640			2.3			 			0.10 U	0.068
Fluorene	86-73-7	320			4.1			 			0.10 U	0.089
Hexachlorobutadiene	87-68-3			2.0 U	1.0 U	2.0 U		 		2.0 U	1.0 U	2.0 U
Hexachlorocyclopentadiene	77-47-4			2.00	1.00	2.00		1		2.00	1.00	2.00
Isophorone	78-59-1							1			1	
· ·	78-59-1 91-20-3		(3)	ana (3)	(4)	(3)		(3)	(3)	(3)	(4)	(4)
Naphthalene		160 (3)	1,200 ⁽³⁾	890 ⁽³⁾	3.6 (4)	120 (3)		0.10 U ⁽³⁾	0.32 (3)	2.0 U ⁽³⁾	0.10 U ⁽⁴⁾	0.020 U ⁽⁴⁾
N-Nitroso-di-n-propylamine	621-64-7	-						-				
N-Nitrosodiphenylamine	86-30-6	-										
Pentachlorophenol	87-86-5	1.0				1					1	1

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Modified from Floyd | Snider 2022 draft RI/FS.

Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

					ndwater Monitoring W	,	11077	NA144 00	ı		B414/ 00	
		Location Name	WCD-MW7-36	WCD-MW07-120511	MW-07 WCD-MW07-112712	WCD-MW07-100113	MW07-WL-03292021	MW-08 WCD-MW8-37	THICK BANKO CE		MW-09 WCD-MW09-112712	WCD-MW09-100113
		Sample Name Sample Date	6/21/2010	12/5/2011	11/27/2012	10/1/2013	3/29/2021	6/22/2010	WCD-MW9-65 6/21/2010	WCD-MW09-120511 12/5/2011	11/27/2012	10/1/2013
			6/21/2010	12/5/2011	11/2//2012	10/1/2013	3/29/2021	6/22/2010	6/21/2010	12/5/2011	11/2//2012	10/1/2013
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds (con	· · · · · · · · · · · · · · · · · · ·					T T	ı	T .	T			
Phenanthrene	85-01-8	-			1.1						0.10 U	0.18
Phenol	108-95-2	4,800										
Pyrene	129-00-0	240			0.20						0.10 U	0.060
Total Naphthalenes		160			5.2						0.10 U	0.023
Volatile Organic Compounds												
1,1,1,2-Tetrachloroethane	630-20-6			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,1,1-Trichloroethane	71-55-6	200		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,1,2,2-Tetrachloroethane	79-34-5			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,1,2-Trichloroethane	79-00-5			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,1-Dichloroethane	75-34-3	7.7		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,1-Dichloroethene	75-35-4			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,1-Dichloropropene	563-58-6			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,2,3-Trichlorobenzene	87-61-6	-		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,2,3-Trichloropropane	96-18-4			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,2,4-Trichlorobenzene	120-82-1			2.0 U	1.0 U	2.0 U		ĺ	Ī	2.0 U	1.0 U	2.0 U
1,2,4-Trimethylbenzene	95-63-6	80		18	1.0 U	2.4				2.0 U	1.0 U	2.0 U
1,2-Dibromo-3-chloropropane	96-12-8			10 U	1.0 U	10 U				10 U	1.0 U	10 U
1,2-Dibromoethane	106-93-4			0.010 U	1.0 U	0.010 U				0.010 U	1.0 U	0.010 U
1,2-Dichlorobenzene	95-50-1			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,2-Dichloroethane	107-06-2	_		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,2-Dichloropropane	78-87-5			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,3,5-Trimethylbenzene	108-67-8	80		6.6	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,3-Dichlorobenzene	541-73-1			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
				2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
1,3-Dichloropropane	142-28-9	-										
1,4-Dichlorobenzene	106-46-7			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
2,2-Dichloropropane	594-20-7	-		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
2,4-Dinitrotoluene	121-14-2	-										
2,6-Dinitrotoluene	606-20-2											
2-Chlorotoluene	95-49-8			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
2-Hexanone	591-78-6			10 U	1.0 U	10 U				10 U	1.0 U	10 U
2-Nitroaniline	88-74-4											
3-Nitroaniline	99-09-2											
4-Chloroaniline	106-47-8											
4-Chlorophenyl phenyl ether	7005-72-3	-										
4-Chlorotoluene	106-43-4			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
4-Nitroaniline	100-01-6											
Acetone	67-64-1			25 U	10 U	25 U				25 U	10 U	25 U
Benzene	71-43-2	5.0		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Bis(2-chloroethyl)ether	111-44-4											
Bis(2-chloroisopropyl)ether	39638-32-9											
Bromobenzene	108-86-1			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Bromochloromethane	74-97-5			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Bromodichloromethane	75-27-4			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Bromoform	75-25-2			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Bromomethane	74-83-9			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Carbon disulfide	75-15-0			2.0 U	=:50	2.0 U		1		2.0 U	-:00	2.0 U
Carbon tetrachloride	56-23-5			2.0 U	1.0 U	2.0 U		1	 	2.0 U	1.0 U	2.0 U
Chlorobenzene	108-90-7			2.0 U	1.0 U	2.0 U		1	 	2.0 U	1.0 U	2.0 U
Chloroethane	75-00-3			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Chloroform	67-66-3	70		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U

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Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (μg/L)

		Location Name			MW-07			MW-08			MW-09	
		Sample Name	WCD-MW7-36	WCD-MW07-120511	WCD-MW07-112712	WCD-MW07-100113	MW07-WL-03292021	WCD-MW8-37	WCD-MW9-65	WCD-MW09-120511	WCD-MW09-112712	WCD-MW09-100113
		Sample Date	6/21/2010	12/5/2011	11/27/2012	10/1/2013	3/29/2021	6/22/2010	6/21/2010	12/5/2011	11/27/2012	10/1/2013
Analyte	CAS No.	Screening Level (1)										
Volatile Organic Compounds (cont.)	•										•	
Chloromethane	74-87-3	-		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
cis-1,2-Dichloroethene	156-59-2	70		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
cis-1,3-Dichloropropene	10061-01-5			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Cymene	99-87-6			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Dibromochloromethane	124-48-1			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Dibromomethane	74-95-3			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Dichlorodifluoromethane	75-71-8			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Ethylbenzene	100-41-4	700		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Hexachlorobenzene	118-74-1											
Hexachloroethane	67-72-1											
Isopropylbenzene	98-82-8	800		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Methyl ethyl ketone	78-93-3			10 U	10 U	10 U				10 U	10 U	10 U
Methyl isobutyl ketone	108-10-1	640		10 U	1.0 U	10 U				10 U	1.0 U	10 U
Methylene chloride	75-09-2			5.0 U	1.0 U	5.0 U				5.0 U	1.0 U	5.0 U
Methyl-tert-butyl ether	1634-04-4			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
n-Butylbenzene	104-51-8			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Nitrobenzene	98-95-3											
n-Propylbenzene	103-65-1	800		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
sec-Butylbenzene	135-98-8			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Styrene	100-42-5	100	1.0 U	2.0 U	1.0 U	2.0 U		1.0 U	1.0 U	2.0 U	1.0 U	2.0 U
tert-Butylbenzene	98-06-6			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Tetrachloroethene	127-18-4			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Toluene	108-88-3	1,000		2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
trans-1,2-Dichloroethene	156-60-5			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
trans-1,3-Dichloropropene	10061-02-6			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Trichloroethene	79-01-6	5.0		2.0 U	1.0 U	2.0 U				2.0 U	6.7	5.9
Trichlorofluoromethane	75-69-4			2.0 U	1.0 U	2.0 U				2.0 U	1.0 U	2.0 U
Vinyl chloride	75-01-4			0.20 U	0.20 U	0.20 U				0.20 U	0.20 U	0.20 U
Xylene (meta & para)	108-38-3/106-42-3			4.0 U		4.0 U				4.0 U		4.0 U
Xylene (ortho)	95-47-6			2.0 U		2.0 U				2.0 U		2.0 U
Xylene (total)	1330-20-7	1,000		4.0 U	1.0 U	4.0 U			_	4.0 U	1.0 U	4.0 U
Other												
PBDE-003	101-55-3											

Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

- Not available or not established. BOLD Detected result exceeds criteria.

The screening level is a proposed Site cleanup level.

1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.

2 A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest RL was selected. 3 Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level for total naphthalenes. 4 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level. 5 Xylene (total) as calculated by Floyd | Snider was preferred over laboratory reported result.

CAS = Chemical Abstracts Service cPAH = Carcinogenic polycyclic aromatic hydrocarbon

PBDE = Polybrominated diphenyl ethers TEQ = Toxic equivalent

μg/L = Micrograms per liter

USEPA = U.S. Environmental Protection Agency

MTCA = Model Toxics Control Act

Qualifiers:

J Analyte was detected and the concentration is estimated.

U Analyte is not detected at the associated reporting limit.

UJ Analyte is not detected at the associated reporting limit, which is an estimate.

Modified from Floyd | Snider 2022 draft RI/FS.



Modified from Floyd | Snider 2022 draft RI/FS.

July 2024 PUBLIC REVIEW DRAFT

Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		MW-09 (cont l		1		N	1W-10		
		Sample Name	MW09-09152020	MW-09-12172020	MW09-032921	MW09-06222021	WCD-MW10-43	WCD-MW10-120611	WCD-MW10-112712	WCD-MW10-030513	WCD-MW10-060413	WCD-MW10-100113
		Sample Name	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.		3/13/2020	12/17/2020	3/23/2021	0/22/2021	0/22/2010	12/0/2011	11/2//2012	3/3/2013	0/4/2013	10/1/2013
	CAS NO.	Screening Level (1)										
Semivolatile Organic Compounds 1-Methylnaphthalene	90-12-0		0.099 U		ı	I	1	I	510		1	
2,4,5-Trichlorophenol	95-95-4	1,600	2.0 U				-		510			
	88-06-2	8.0	2.0 U				-					
2,4,6-Trichlorophenol 2,4-Dichlorophenol	120-83-2	48	2.0 U									
		1	0.99 U									
2,4-Dimethylphenol	105-67-9 51-28-5	320 32	0.99 U 2.0 UJ									
2,4-Dinitrophenol			2.0 01									
2-Chloronaphthalene	91-58-7		0.00.11									
2-Chlorophenol	95-57-8	40	0.99 U									
2-Methylnaphthalene	91-57-6	-	0.099 U						430			
2-Methylphenol	95-48-7	800	0.99 U									
2-Nitrophenol	88-75-5	-	2.0 U									
3- & 4-Methylphenol	15831-10-4	800	0.99 U									
4,6-Dinitro-o-cresol	534-52-1	1.3	5.0 UJ									
4-Chloro-3-methylphenol	59-50-7	1,600	5.0 U									
4-Nitrophenol	100-02-7	1,600	5.0 UJ									
Acenaphthene	83-32-9	480	0.20 U						6.9			
Acenaphthylene	208-96-8		0.20 U						240			
Acrylonitrile	107-13-1							10 U		10 U	10 U	10 U
Anthracene	120-12-7	2,400	0.099 U						0.10 U			
Benzo(a)anthracene	56-55-3		0.20 U				0.10 U		0.10 U			
Benzo(a)pyrene	50-32-8		0.099 U				0.10 U		0.10 U			
Benzo(b)fluoranthene	205-99-2		0.20 U				0.10 U		0.10 U			
Benzo(g,h,i)perylene	191-24-2		0.20 U						0.10 U			
Benzo(k)fluoranthene	207-08-9		0.20 U				0.10 U		0.10 U			
Benzoic acid	65-85-0											
Benzyl alcohol	100-51-6											
Bis(2-chloroethoxy)methane	111-91-1											
Bis(2-ethylhexyl)phthalate	117-81-7	-										
Butyl benzyl phthalate	85-68-7	-										
Carbazole	86-74-8	-										
Chrysene	218-01-9	-	0.20 U				0.10 U		0.10 U			
Dibenzo(a,h)anthracene	53-70-3		0.20 U				0.10 U		0.10 U			
Indeno(1,2,3-c,d)pyrene	193-39-5		0.099 U				0.10 U		0.10 U			
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	0.099 U				0.10 U		0.10 U			
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	0.099 U				0.10 U		0.10 U			
Dibenzofuran	132-64-9											
Diethylphthalate	84-66-2						1					
Dimethyl phthalate	131-11-3											
Di-n-butyl phthalate	84-74-2						1				İ	
Di-n-octyl phthalate	117-84-0						İ				1	
Fluoranthene	206-44-0	640	0.099 U				İ		0.10 U		1	
Fluorene	86-73-7	320	0.099 U				İ		0.10 U		1	
Hexachlorobutadiene	87-68-3						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Hexachlorocyclopentadiene	77-47-4											
Isophorone	78-59-1						!					
Naphthalene	160 ⁽³⁾	160 ⁽³⁾	2.0 U ⁽³⁾	1.0 U ⁽³⁾	1.3 U ⁽³⁾	1 2 (3)	10,000 ⁽³⁾	14,000 ⁽³⁾	11,000 (4)	2,300 (3)	12 000 (3)	12,000 ⁽³⁾
N-Nitroso-di-n-propylamine	621-64-7	160 ***	2.0 0 * *	1.00	1.3 U	1.3 U ⁽³⁾	10,000	14,000	11,000	2,300	12,000 ⁽³⁾	12,000
N-Nitrosodiphenylamine	86-30-6										1	
Pentachlorophenol	87-86-5	1.0	0.099 UJ		-		 				-	
i ciitaciiioropiielloi	07-00-3	±.∪	U.U55 UJ		1	l		1	l		1	l .



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		MW-09 (cont)		1		Ι.	/W-10		
		Sample Name	MW09-09152020	MW-09-12172020	MW09-032921	MW09-06222021	WCD-MW10-43	WCD-MW10-120611	WCD-MW10-112712	WCD-MW10-030513	WCD-MW10-060413	WCD-MW10-100113
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.		3/13/2020	12/17/2020	3/23/2021	0/22/2021	0/22/2010	12/0/2011	11/2//2012	3/3/2013	0/4/2013	10/1/2013
Semivolatile Organic Compounds (con		Screening Level (1)										
Phenanthrene	85-01-8		0.099 U		1		ı		0.10 U			
Phenol	108-95-2	4,800	2.0 UJ						0.10 0			
Pyrene	129-00-0	240	0.20 U						0.10 U			
Total Naphthalenes		160	2.0 U						12,000			
Volatile Organic Compounds		100	2.0 0						12,000			
1,1,1,2-Tetrachloroethane	630-20-6	- 1			1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1.1.1-Trichloroethane	71-55-6	200	1.0 U	1.0 U	0.40 U	0.40 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,2,2-Tetrachloroethane	79-34-5		1.00	1.00	0.40 0	0.40 0	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,2-Trichloroethane	79-00-5	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloroethane	75-34-3	7.7	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloroethene	75-35-4		1.00	1.00	0.50 0	0.30 0	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
	563-58-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloropropene 1,2,3-Trichlorobenzene	87-61-6				+		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,3-Trichloropenzene	96-18-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
	120-82-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,4-Trichlorobenzene	95-63-6	 80	1011	10.11	0.50.11	0.50 U	230	310	440	2.0 0	2.0 0 170	2.0 O
1,2,4-Trimethylbenzene 1,2-Dibromo-3-chloropropane			1.0 U	1.0 U	0.50 U	0.50 0	10 U	10 U	1.0 U	10 U	170 10 U	150 10 U
	96-12-8 106-93-4				-		1.0 U	0.010 U	1.0 U	0.010 U	0.010 U	0.010 U
1,2-Dibromoethane 1,2-Dichlorobenzene					-		1.0 U	0.010 U	1.0 U	2.0 U	2.0 U	2.0 U
	95-50-1 107-06-2				-		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2-Dichloroethane					-		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2-Dichloropropane	78-87-5 108-67-8	 80	40.11	40.11	0.25.11	0.25 U	90	2.0 0 110	87	2.0 0	63	2.0 0
1,3,5-Trimethylbenzene			1.0 U	1.0 U	0.25 U	0.25 0		2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,3-Dichlorobenzene	541-73-1				-		1.0 U	2.0 U		2.0 U		
1,3-Dichloropropane	142-28-9 106-46-7				-		1.0 U 1.0 U	2.0 U	1.0 U 1.0 U	2.0 U	2.0 U 2.0 U	2.0 U 2.0 U
1,4-Dichlorobenzene	594-20-7						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2,2-Dichloropropane					-		1.00	2.0 0	1.00	2.0 0	2.0 0	2.0 0
2,4-Dinitrotoluene	121-14-2				-							
2,6-Dinitrotoluene	606-20-2				-		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2-Chlorotoluene	95-49-8				-							
2-Hexanone	591-78-6 88-74-4				-		10 U	10 U	1.0 U	10 U	10 U	10 U
2-Nitroaniline					-							
3-Nitroaniline	99-09-2 106-47-8				-							
4-Chloroaniline	7005-72-3				-							
4-Chlorophenyl phenyl ether					-		1011	2011	1011	2011	2011	2011
4-Chlorotoluene	106-43-4				 		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
4-Nitroaniline	100-01-6 67-64-1				 		10 U	25 U	10 U	25 U	25 U	25 U
Acetone			1.0 U	1.0 U	0.44 U	0.44 U	10 U 1.3	25 U 2.0 U	10 U 1.6	25 U 2.0 U	25 U 2.0 U	25 U 2.0 U
Benzene	71-43-2	5.0	1.00	1.00	0.44 U	U.44 U	1.3	2.0 0	1.0	2.0 0	2.0 0	2.0 0
Bis(2-chloroethyl)ether	111-44-4				 					 		
Bis(2-chloroisopropyl)ether	39638-32-9				.		16	2011	10	2011	2011	2011
Bromobenzene	108-86-1				!		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromochloromethane	74-97-5				-		40	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromodichloromethane	75-27-4				-		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromoform	75-25-2						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromomethane	74-83-9				ļ		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Carbon disulfide	75-15-0							2.0 U		2.0 U	2.0 U	2.0 U
Carbon tetrachloride	56-23-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chlorobenzene	108-90-7						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chloroethane	75-00-3	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chloroform	67-66-3	70	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U



Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (μg/L)

		Location Name		MW-09 (cont.)				N	1W-10		
		Sample Name	MW09-09152020	MW-09-12172020	MW09-032921	MW09-06222021	WCD-MW10-43	WCD-MW10-120611	WCD-MW10-112712	WCD-MW10-030513	WCD-MW10-060413	WCD-MW10-100113
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.	Screening Level (1)										
Volatile Organic Compounds (cont.)												
Chloromethane	74-87-3		2.0 U	2.0 U	0.75 U	0.75 U	10 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
cis-1,2-Dichloroethene	156-59-2	70	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
cis-1,3-Dichloropropene	10061-01-5	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Cymene	99-87-6						11	13	9.8	29	22	24
Dibromochloromethane	124-48-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Dibromomethane	74-95-3						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Dichlorodifluoromethane	75-71-8						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Ethylbenzene	100-41-4	700	1.0 U	1.0 U	0.40 U	0.76	660	720	890	520	330	360
Hexachlorobenzene	118-74-1											
Hexachloroethane	67-72-1											
Isopropylbenzene	98-82-8	800	1.0 U	1.0 U	0.50 U	0.50 U	41	42	40	36	27	28
Methyl ethyl ketone	78-93-3	-			1.5 U		10 U	10 U	10 U	10 U	10 U	10 U
Methyl isobutyl ketone	108-10-1	640	1.3 U	1.3 U		1.3 U	10 U	10 U	1.0 U	10 U	10 U	10 U
Methylene chloride	75-09-2	-					5.0 U	5.0 U	1.0 U	5.0 U	5.0 U	5.0 U
Methyl-tert-butyl ether	1634-04-4	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
n-Butylbenzene	104-51-8	-						2.0 U	6.4	2.0 U	2.0 U	2.0 U
Nitrobenzene	98-95-3	-										
n-Propylbenzene	103-65-1	800	1.0 U	1.0 U	0.50 U	0.50 U	14	15	16	2.0 U	12	12
sec-Butylbenzene	135-98-8	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Styrene	100-42-5	100	1.0 U	1.0 U	0.50 U	0.50 U	15	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
tert-Butylbenzene	98-06-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Tetrachloroethene	127-18-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Toluene	108-88-3	1,000	1.0 U	1.0 U	0.75 U	0.75 U	61	71	52	49	40	40
trans-1,2-Dichloroethene	156-60-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
trans-1,3-Dichloropropene	10061-02-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Trichloroethene	79-01-6	5.0	2.7	2.5	1.7	2.1	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Trichlorofluoromethane	75-69-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Vinyl chloride	75-01-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U
Xylene (meta & para)	108-38-3/106-42-3		1.0 U	1.0 U	1.0 U	1.0 U	790	950		660	490	500
Xylene (ortho)	95-47-6		1.0 U	1.0 U	0.50 U	0.50 U	460	510		370	310	290
Xylene (total)	1330-20-7	1,000	1.0 U ⁽⁵⁾	1.0 U	1.0 U ⁽⁵⁾	1.0 U ⁽⁵⁾	1,200	1,500	1,900	1,000	800	790
Other												
PBDE-003	101-55-3											

Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

- Not available or not established.

BOLD Detected result exceeds criteria.

- 1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landay Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.
- 2 A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest RL was selected. 3 Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level for total naphthalenes.
- 4 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level. 5 Xylene (total) as calculated by Floyd | Snider was preferred over laboratory reported result.

Abbreviations:

CAS = Chemical Abstracts Service

PBDE = Polybrominated diphenyl ethers

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

 μ g/L = Micrograms per liter MTCA = Model Toxics Control Act USEPA = U.S. Environmental Protection Agency

- J Analyte was detected and the concentration is estimated.
- U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table~5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (µg/L)

						ng wen Analytica		11-07-7				
		Location Name	BAN440 004 F 2020	MW-10 (PWM10 00333031	14/CD 8414/44 42	MCD 141444 420544		1W-11	14/CD 14/4/14 0C0443	WCD 141444 400443
		Sample Name	MW10-09152020	MW-10-12172020		MW10-06222021	WCD-MW11-42			WCD-MW11-030513		
Analyte	CAS No.	Sample Date Screening Level (1)	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/5/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Semivolatile Organic Compounds		Screening Level					<u> </u>				1	
1-Methylnaphthalene	90-12-0								0.10 U			
2,4,5-Trichlorophenol	95-95-4	1,600										
2,4,6-Trichlorophenol	88-06-2	8.0										
2,4-Dichlorophenol	120-83-2	48										
2,4-Dimethylphenol	105-67-9	320										
2,4-Dinitrophenol	51-28-5	32										
2-Chloronaphthalene	91-58-7											
2-Chlorophenol	95-57-8	40										
2-Methylnaphthalene	91-57-6								0.10 U			
2-Methylphenol	95-48-7	800					1					
2-Nitrophenol	88-75-5						1					
3- & 4-Methylphenol	15831-10-4	800										
4,6-Dinitro-o-cresol	534-52-1	1.3					1					
4-Chloro-3-methylphenol	59-50-7	1,600					1					
4-Nitrophenol	100-02-7	1,600					1					
Acenaphthene	83-32-9	480							0.10 U			
Acenaphthylene	208-96-8								0.10 U			
Acrylonitrile	107-13-1							10 U	0.20 0	10 U	10 U	10 U
Anthracene	120-12-7	2,400						10 0	0.10 U	10 0	10 0	10 0
Benzo(a)anthracene	56-55-3	2,400					0.10 U		0.10 U			
	50-32-8	-					0.10 U		0.10 U			
Benzo(a)pyrene	205-99-2						0.10 U		0.10 U			
Benzo(b)fluoranthene	191-24-2						0.10 0		0.10 U			
Benzo(g,h,i)perylene Benzo(k)fluoranthene	207-08-9						0.10 U		0.10 U			
	65-85-0						0.10 0		0.10 0			
Benzoic acid												
Benzyl alcohol	100-51-6											
Bis(2-chloroethoxy)methane	111-91-1											
Bis(2-ethylhexyl)phthalate	117-81-7	-										
Butyl benzyl phthalate	85-68-7											
Carbazole	86-74-8	-					0.4011		0.40.11			
Chrysene	218-01-9						0.10 U		0.10 U			
Dibenzo(a,h)anthracene	53-70-3						0.10 U		0.10 U			
Indeno(1,2,3-c,d)pyrene	193-39-5						0.10 U		0.10 U			
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10					0.10 U		0.10 U			
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10			1		0.10 U		0.10 U		1	
Dibenzofuran	132-64-9						1					
Diethylphthalate	84-66-2						!				ļ	
Dimethyl phthalate	131-11-3				1		 				1	
Di-n-butyl phthalate	84-74-2											
Di-n-octyl phthalate	117-84-0						1					
Fluoranthene	206-44-0	640					ļ		0.10 U			
Fluorene	86-73-7	320					ļ		0.10 U			
Hexachlorobutadiene	87-68-3	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Hexachlorocyclopentadiene	77-47-4	-										
Isophorone	78-59-1											
Naphthalene	160 ⁽³⁾	160 ⁽³⁾	7,700 ⁽³⁾	17,000 ⁽³⁾	11,000 ⁽³⁾	16,000 ⁽³⁾	3.1 (3)	2.1 (3)	0.10 U ⁽⁴⁾	380 ⁽³⁾	2.0 U ⁽³⁾	7.6 (3)
N-Nitroso-di-n-propylamine	621-64-7											
N-Nitrosodiphenylamine	86-30-6											
Pentachlorophenol	87-86-5	1.0										

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Modified from Floyd | Snider 2022 draft RI/FS.

Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		MW-10 (c	ont.)				N	MW-11		-
		Sample Name	MW10-09152020	MW-10-12172020	MW10-032921	MW10-06222021	WCD-MW11-42	WCD-MW11-120511	WCD-MW11-112712	WCD-MW11-030513	WCD-MW11-060413	WCD-MW11-100113
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/5/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.	Screening Level (1)	3/13/2020	12/1//2020	0/25/2021	0/22/2021	0/22/2010	12/5/2011	11/2//2012	5/5/2015	07-172020	10/1/2010
Semivolatile Organic Compounds (co		Screening Level										
				ı		1		1	0.4011	1		
Phenanthrene	85-01-8								0.10 U			
Phenol	108-95-2	4,800							-			
Pyrene	129-00-0	240							0.10 U			
Total Naphthalenes		160					<u> </u>	l	0.10 U			
				•	Vo	latile Organic Compou	_					
1,1,1,2-Tetrachloroethane	630-20-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,1-Trichloroethane	71-55-6	200	1.0 U	10 U	8.0 U	8.0 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,2,2-Tetrachloroethane	79-34-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,2-Trichloroethane	79-00-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloroethane	75-34-3	7.7	1.0 U	10 U	10 U	10 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloroethene	75-35-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloropropene	563-58-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,3-Trichlorobenzene	87-61-6	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,3-Trichloropropane	96-18-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,4-Trichlorobenzene	120-82-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,4-Trimethylbenzene	95-63-6	80	260	330	250	200	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2-Dibromo-3-chloropropane	96-12-8						10 U	10 U	1.0 U	10 U	10 U	10 U
1,2-Dibromoethane	106-93-4	-					1.0 U	0.010 U	1.0 U	0.010 U	0.010 U	0.010 U
1,2-Dichlorobenzene	95-50-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2-Dichloroethane	107-06-2	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2-Dichloropropane	78-87-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1.3.5-Trimethylbenzene	108-67-8	80	75	110	87	62	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,3-Dichlorobenzene	541-73-1		,,,			V-	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,3-Dichloropropane	142-28-9						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,4-Dichlorobenzene	106-46-7						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2,2-Dichloropropane	594-20-7	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2.4-Dinitrotoluene	121-14-2						1.00	2.0 0	1.0 0	2.0 0	2.00	2.0 0
2.6-Dinitrotoluene	606-20-2								+			
2-Chlorotoluene	95-49-8						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2-Hexanone	591-78-6	-					1.0 U	2.0 U	1.0 U	2.0 U	10 U	2.0 U
2-Hexanone 2-Nitroaniline	88-74-4						10 0	10 0	1.00	10 0	10 0	10 0
		-										—
3-Nitroaniline 4-Chloroaniline	99-09-2											
	106-47-8	-										—
4-Chlorophenyl phenyl ether	7005-72-3											
4-Chlorotoluene	106-43-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
4-Nitroaniline	100-01-6											
Acetone	67-64-1						10 U	25 U	10 U	25 U	25 U	25 U
Benzene	71-43-2	5.0	1.0 U	10 U	8.8 U	8.8 U	0.35 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bis(2-chloroethyl)ether	111-44-4											
Bis(2-chloroisopropyl)ether	39638-32-9											
Bromobenzene	108-86-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromochloromethane	74-97-5							2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromodichloromethane	75-27-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromoform	75-25-2						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromomethane	74-83-9	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Carbon disulfide	75-15-0							2.0 U		2.0 U	2.0 U	2.0 U
Carbon tetrachloride	56-23-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chlorobenzene	108-90-7						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chloroethane	75-00-3						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chloroform	67-66-3	70	1.0 U	10 U	10 U	10 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U

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Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (µg/L)

						ing wen Analytica		- (1-0) -/				
		Location Name		MW-10 (/W-11		
		Sample Name	MW10-09152020	MW-10-12172020	MW10-032921	MW10-06222021	WCD-MW11-42	WCD-MW11-120511	WCD-MW11-112712	WCD-MW11-030513	WCD-MW11-060413	WCD-MW11-100113
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/5/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.	Screening Level (1)										
Volatile Organic Compounds (cont.	.)											
Chloromethane	74-87-3	-	2.0 U	20 U	15 U	15 U	10 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
cis-1,2-Dichloroethene	156-59-2	70	1.0 U	10 U	10 U	10 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
cis-1,3-Dichloropropene	10061-01-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Cymene	99-87-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Dibromochloromethane	124-48-1	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Dibromomethane	74-95-3	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Dichlorodifluoromethane	75-71-8	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Ethylbenzene	100-41-4	700	570	690	540	470	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Hexachlorobenzene	118-74-1	-										
Hexachloroethane	67-72-1											
Isopropylbenzene	98-82-8	800	32	35	27	22	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Methyl ethyl ketone	78-93-3				30 U		10 U	10 U	10 U	10 U	10 U	10 U
Methyl isobutyl ketone	108-10-1	640	1.3 U	13 U		25 U	10 U	10 U	1.0 U	10 U	10 U	10 U
Methylene chloride	75-09-2						5.0 U	5.0 U	1.0 U	5.0 U	5.0 U	5.0 U
Methyl-tert-butyl ether	1634-04-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
n-Butylbenzene	104-51-8							2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Nitrobenzene	98-95-3	-										
n-Propylbenzene	103-65-1	800	12	13	10 U	10 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
sec-Butylbenzene	135-98-8						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Styrene	100-42-5	100	1.0 U	10 U	10 U	10 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
tert-Butylbenzene	98-06-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Tetrachloroethene	127-18-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Toluene	108-88-3	1,000	38	38	31	34	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
trans-1,2-Dichloroethene	156-60-5	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
trans-1,3-Dichloropropene	10061-02-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Trichloroethene	79-01-6	5.0	0.50 U	5.0 U	10 U	10 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Trichlorofluoromethane	75-69-4	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Vinyl chloride	75-01-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U
Xylene (meta & para)	108-38-3/106-42-3		730	850	760	590	2.0 U	4.0 U		4.0 U	4.0 U	4.0 U
Xylene (ortho)	95-47-6		420	450	410	360	1.0 U	2.0 U		2.0 U	2.0 U	2.0 U
Xylene (total)	1330-20-7	1,000	1,200 (5)	1,300	1,200 ⁽⁵⁾	940 ⁽⁵⁾	2.0 U	4.0 U	1.0 U	4.0 U	4.0 U	4.0 U
Other	•			•			•		•			
PBDE-003	101-55-3											
					•			1				1

Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

- Not available or not established.

BOLD Detected result exceeds criteria.

The screening level is a proposed Site cleanup level.

- 1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial investigation and Feasibility Study apart from updating screening levels.
- 2.A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest RL was selected. 3. Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level for total naphthalenes. 4 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level. 5 Xylene (total) as calculated by Floyd | Snider was preferred over laboratory reported result.

Abbreviations:

CAS = Chemical Abstracts Service

PBDE = Polybrominated diphenyl ethers TEQ = Toxic equivalent

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

μg/L = Micrograms per liter MTCA = Model Toxics Control Act USEPA = U.S. Environmental Protection Agency

Qualifiers:

J Analyte was detected and the concentration is estimated.

U Analyte is not detected at the associated reporting limit.

Modified from Floyd | Snider 2022 draft RI/FS.

July 2024 PUBLIC REVIEW DRAFT

UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		MW-11	(cont.)					MW-12		
		Sample Name	MW11-09152020	MW-11-12172020	MW11-032921	MW11-06222021	WCD-MW12-42	WCD-MW12-120611	WCD-MW12-112712	WCD-MW12-030513	WCD-MW12-060413	WCD-MW12-100113
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds												
1-Methylnaphthalene	90-12-0	-							0.10 U			
2,4,5-Trichlorophenol	95-95-4	1,600										
2,4,6-Trichlorophenol	88-06-2	8.0										
2,4-Dichlorophenol	120-83-2	48										
2,4-Dimethylphenol	105-67-9	320										
2,4-Dinitrophenol	51-28-5	32										
2-Chloronaphthalene	91-58-7											
2-Chlorophenol	95-57-8	40										
2-Methylnaphthalene	91-57-6								0.10 U			
2-Methylphenol	95-48-7	800										
2-Nitrophenol	88-75-5											
3- & 4-Methylphenol	15831-10-4	800										
4,6-Dinitro-o-cresol	534-52-1	1.3										
4-Chloro-3-methylphenol	59-50-7	1,600										
4-Nitrophenol	100-02-7	1,600										
Acenaphthene	83-32-9	480							2.7			
Acenaphthylene	208-96-8								7.8			
Acrylonitrile	107-13-1							10 U		10 U	10 U	10 U
Anthracene	120-12-7	2,400							0.10 U			
Benzo(a)anthracene	56-55-3						0.27		0.10 U			
Benzo(a)pyrene	50-32-8						0.10 U		0.10 U			
Benzo(b)fluoranthene	205-99-2						0.10 U		0.10 U			
Benzo(g,h,i)perylene	191-24-2								0.10 U			
Benzo(k)fluoranthene	207-08-9						0.10 U		0.10 U			
Benzoic acid	65-85-0											
Benzyl alcohol	100-51-6											
Bis(2-chloroethoxy)methane	111-91-1											
Bis(2-ethylhexyl)phthalate	117-81-7											
Butyl benzyl phthalate	85-68-7											
Carbazole	86-74-8											
Chrysene	218-01-9						0.23		0.10 U			
Dibenzo(a,h)anthracene	53-70-3						0.10 U		0.10 U			
Indeno(1,2,3-c,d)pyrene	193-39-5						0.10 U		0.10 U			
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10					0.099		0.10 U			
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10					0.029		0.10 U			
Dibenzofuran	132-64-9											
Diethylphthalate	84-66-2											
Dimethyl phthalate	131-11-3								İ			İ
Di-n-butyl phthalate	84-74-2											
Di-n-octyl phthalate	117-84-0								1			1
Fluoranthene	206-44-0	640							1.8			İ
Fluorene	86-73-7	320							0.10 U			1
Hexachlorobutadiene	87-68-3						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Hexachlorocyclopentadiene	77-47-4				1	1			T			T
Isophorone	78-59-1											
Naphthalene	91-20-3	160 ⁽³⁾	220 (3)	15 ⁽³⁾	1.3 U ⁽³⁾	1.3 U ⁽³⁾	8,400 ⁽³⁾	11,000 ⁽³⁾	0.10 U ⁽⁴⁾	210 (3)	12,000 ⁽³⁾	7,500 ⁽³⁾
N-Nitroso-di-n-propylamine	621-64-7		220	13	1.3 0	1.5 0	0,400	11,000	0.10 0	210	12,000	7,300
N-Nitrosodiphenylamine	86-30-6				1	1	1		 			
Pentachlorophenol	87-86-5	1.0					 		 			
i cittacinorophenor	07-00-3	1.0	l		l	1						

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Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		MW-11	(cont.)					MW-12		
		Sample Name	MW11-09152020	MW-11-12172020	MW11-032921	MW11-06222021	WCD-MW12-42	WCD-MW12-120611	WCD-MW12-112712	WCD-MW12-030513	WCD-MW12-060413	WCD-MW12-100113
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds (co	ont.)	Screening Level										
Phenanthrene	85-01-8								0.10 U			
Phenol	108-95-2	4,800										
Pyrene	129-00-0	240							1.3			
Total Naphthalenes		160							0.10 U			
Volatile Organic Compounds	•											
1,1,1,2-Tetrachloroethane	630-20-6	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,1-Trichloroethane	71-55-6	200	1.0 U	1.0 U	0.40 U	0.40 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,2,2-Tetrachloroethane	79-34-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1,2-Trichloroethane	79-00-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloroethane	75-34-3	7.7	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloroethene	75-35-4	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,1-Dichloropropene	563-58-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,3-Trichlorobenzene	87-61-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,3-Trichloropropane	96-18-4	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,4-Trichlorobenzene	120-82-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2,4-Trimethylbenzene	95-63-6	80	1.0 U	1.0 U	0.50 U	0.50 U	270	300	1.1	13	230	230
1,2-Dibromo-3-chloropropane	96-12-8						10 U	10 U	1.0 U	10 U	10 U	10 U
1,2-Dibromoethane	106-93-4						1.0 U	0.010 U	1.0 U	0.010 U	0.010 U	0.010 U
1,2-Dichlorobenzene	95-50-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2-Dichloroethane	107-06-2	_					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,2-Dichloropropane	78-87-5	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,3,5-Trimethylbenzene	108-67-8	80	1.0 U	1.0 U	0.25 U	0.25 U	110	110	1.0 U	3.5	86	96
1,3-Dichlorobenzene	541-73-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,3-Dichloropropane	142-28-9						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
1,4-Dichlorobenzene	106-46-7	-					1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2,2-Dichloropropane	594-20-7						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2,4-Dinitrotoluene	121-14-2											
2,6-Dinitrotoluene	606-20-2											
2-Chlorotoluene	95-49-8						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
2-Hexanone	591-78-6						10 U	10 U	1.0 U	10 U	10 U	10 U
2-Nitroaniline	88-74-4	_					10 0	10 0	1.0 0	10 0	10 0	10 0
3-Nitroaniline	99-09-2											
4-Chloroaniline	106-47-8	_										
4-Chlorophenyl phenyl ether	7005-72-3											
4-Chlorotoluene	106-43-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
4-Nitroaniline	100-01-6											
Acetone	67-64-1	_					10 U	25 U	10 U	25 U	25 U	25 U
Benzene	71-43-2	5.0	1.0 U	1.0 U	0.44 U	0.44 U	13	12	1.0 U	2.0 U	14	13
Bis(2-chloroethyl)ether	111-44-4				*****							
Bis(2-chloroisopropyl)ether	39638-32-9											
Bromobenzene	108-86-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromochloromethane	74-97-5						1.0 0	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromodichloromethane	75-27-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromoform	75-25-2				<u> </u>		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Bromomethane	74-83-9	-			<u> </u>		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Carbon disulfide	75-15-0						1.0 0	2.0 U	1.00	2.0 U	2.0 U	2.0 U
Carbon tetrachloride	56-23-5				1		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chlorobenzene	108-90-7				+		1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chloroethane	75-00-3						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Chloroform	67-66-3	70	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U



Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (µg/L)

		Location Name		MW-11	(cont.)					MW-12		
		Sample Name	MW11-09152020	MW-11-12172020	MW11-032921	MW11-06222021	WCD-MW12-42	WCD-MW12-120611	WCD-MW12-112712	WCD-MW12-030513	WCD-MW12-060413	WCD-MW12-10011
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	6/22/2010	12/6/2011	11/27/2012	3/5/2013	6/4/2013	10/1/2013
Analyte	CAS No.	Screening Level (1)	-,,		-,,	-,,	-,,	, _,		0,0,000	-, -,	
Volatile Organic Compounds (cont		bercenning zever										
Chloromethane	74-87-3		2.0 U	2.0 U	0.75 U	0.75 U	10 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
cis-1.2-Dichloroethene	156-59-2	70	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
cis-1,3-Dichloropropene	10061-01-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Cymene	99-87-6						10	13	1.0 U	2.7	53	39
Dibromochloromethane	124-48-1						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Dibromomethane	74-95-3						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Dichlorodifluoromethane	75-71-8						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Ethylbenzene	100-41-4	700	1.0 U	1.0 U	0.40 U	0.40 U	1,000	890	1.4	38	1,000	900
Hexachlorobenzene	118-74-1											
Hexachloroethane	67-72-1											
Isopropylbenzene	98-82-8	800	1.0 U	1.0 U	0.50 U	0.50 U	55	46	1.0 U	2.2	40	37
Methyl ethyl ketone	78-93-3				1.5 U		10 U	10 U	10 U	10 U	10 U	10 U
Methyl isobutyl ketone	108-10-1	640	1.3 U	1.3 U		1.3 U	10 U	10 U	1.0 U	10 U	10 U	10 U
Methylene chloride	75-09-2						5.0 U	5.0 U	1.0 U	5.0 U	5.0 U	5.0 U
Methyl-tert-butyl ether	1634-04-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
n-Butylbenzene	104-51-8							2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Nitrobenzene	98-95-3											
n-Propylbenzene	103-65-1	800	1.0 U	1.0 U	0.50 U	0.50 U	16	16	1.0 U	2.0 U	17	15
sec-Butylbenzene	135-98-8						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Styrene	100-42-5	100	1.0 U	1.0 U	0.50 U	0.50 U	1.2	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
tert-Butvlbenzene	98-06-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Tetrachloroethene	127-18-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Toluene	108-88-3	1,000	1.0 U	1.0 U	0.75 U	0.75 U	270	350	1.0 U	15	370	370
trans-1,2-Dichloroethene	156-60-5						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
trans-1,3-Dichloropropene	10061-02-6						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Trichloroethene	79-01-6	5.0	0.50 U	0.50 U	0.50 U	0.50 U	1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Trichlorofluoromethane	75-69-4						1.0 U	2.0 U	1.0 U	2.0 U	2.0 U	2.0 U
Vinyl chloride	75-01-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U
Xylene (meta & para)	108-38-3/106-42-3		1.0 U	1.0 U	1.0 U	1.0 U	1,100	940		16	1,100	1,000
Xylene (ortho)	95-47-6		1.0 U	1.0 U	0.50 U	0.50 U	600	530		30	610	530
Xylene (total)	1330-20-7	1,000	1.0 U ⁽⁵⁾	1.0 U	1.0 U (5)	1.0 U (5)	1,700	1,500	1.0 U	46	1,700	1,500
Other	•			•	•	•			•			•
PBDE-003	101-55-3											

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Notes:

Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

- Not available or not established. BOLD Detected result exceeds criteria.

The screening level is a proposed Site cleanup level.

1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.

2 A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest RL was selected. 3 Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level for total naphthalenes.

4 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level. 5 Xylene (total) as calculated by Floyd | Snider was preferred over laboratory reported result.

Abbreviations:

CAS = Chemical Abstracts Service

PBDE = Polybrominated diphenyl ethers TEQ = Toxic equivalent

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

USEPA = U.S. Environmental Protection Agency

μg/L = Micrograms per liter MTCA = Model Toxics Control Act

Qualifiers:

J Analyte was detected and the concentration is estimated.

U Analyte is not detected at the associated reporting limit.

Modified from Floyd | Snider 2022 draft RI/FS.

UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		MW-12	(cont.)			MW-	-14		MW	/-15
		Sample Name	MW12-09152020	MW-12-12172020	MW12-032921	MW12-06222021	MW14-09162020	MW-14-12172020	MW14-032921	MW14-06222021	MW15-09162020	MW-15-12172020
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	9/16/2020	12/17/2020	3/29/2021	6/22/2021	9/16/2020	12/17/2020
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds												
1-Methylnaphthalene	90-12-0						310				0.098 U	
2,4,5-Trichlorophenol	95-95-4	1,600					2.0 U				2.0 U	
2,4,6-Trichlorophenol	88-06-2	8.0					2.0 U				2.0 U	
2,4-Dichlorophenol	120-83-2	48					2.0 U				2.0 U	
2,4-Dimethylphenol	105-67-9	320					1.0				0.98 U	
2,4-Dinitrophenol	51-28-5	32					2.0 UJ				2.0 UJ	
2-Chloronaphthalene	91-58-7	-										
2-Chlorophenol	95-57-8	40					1.0 U				0.98 U	
2-Methylnaphthalene	91-57-6						240				0.098 U	
2-Methylphenol	95-48-7	800					1.0 U				0.98 U	
2-Nitrophenol	88-75-5						2.0 U				2.0 U	
3- & 4-Methylphenol	15831-10-4	800					1.9				0.98 U	
4,6-Dinitro-o-cresol	534-52-1	1.3					500 U				4.9 UJ	
4-Chloro-3-methylphenol	59-50-7	1,600					5.0 U				4.9 U	
4-Nitrophenol	100-02-7	1,600					5.0 UJ				4.9 UJ	
Acenaphthene	83-32-9	480					160				0.20 U	
Acenaphthylene	208-96-8						22				0.20 U	
Acrylonitrile	107-13-1	-									0.20 0	
Anthracene	120-12-7	2,400					7.6				0.098 U	
Benzo(a)anthracene	56-55-3						0.20 U				0.20 U	
Benzo(a)pyrene	50-32-8						0.10 U				0.098 U	
Benzo(b)fluoranthene	205-99-2	-					0.20 U				0.20 U	
Benzo(g,h,i)perylene	191-24-2						0.20 U				0.20 U	
Benzo(k)fluoranthene	207-08-9						0.20 U				0.20 U	
Benzoic acid	65-85-0	-					0.20 0				0.20 0	
Benzyl alcohol	100-51-6											
Bis(2-chloroethoxy)methane	111-91-1	_										
Bis(2-ethylhexyl)phthalate	117-81-7											
Butyl benzyl phthalate	85-68-7											
	86-74-8											
Carbazole							0.20 U				0.20 U	
Chrysene	218-01-9											
Dibenzo(a,h)anthracene	53-70-3	-					0.20 U				0.20 U	
Indeno(1,2,3-c,d)pyrene	193-39-5						0.10 U				0.098 U	
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10					0.10 U				0.098 U	
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10					0.10 U				0.098 U	
Dibenzofuran	132-64-9			 								
Diethylphthalate	84-66-2											
Dimethyl phthalate	131-11-3											
Di-n-butyl phthalate	84-74-2											
Di-n-octyl phthalate	117-84-0	-										
Fluoranthene	206-44-0	640					2.2				0.098 U	
Fluorene	86-73-7	320					58				0.098 U	
Hexachlorobutadiene	87-68-3											
Hexachlorocyclopentadiene	77-47-4											
Isophorone	78-59-1											
Naphthalene	91-20-3	160 ⁽³⁾	4,800 ⁽³⁾	14,000 ⁽³⁾	13,000 ⁽³⁾	17,000 ⁽³⁾	6,700 ⁽³⁾	3,600 ⁽³⁾	5,300 ⁽³⁾	7,900 ⁽³⁾	2.0 U ⁽³⁾	1.0 U ⁽³⁾
N-Nitroso-di-n-propylamine	621-64-7											
N-Nitrosodiphenylamine	86-30-6											
Pentachlorophenol	87-86-5	1.0					0.10 UJ				0.098 UJ	



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

		Location Name		MW-12	(cont.)			MW	-14		MW	<i>l</i> -15
		Sample Name	MW12-09152020	MW-12-12172020	MW12-032921	MW12-06222021	MW14-09162020	MW-14-12172020	MW14-032921	MW14-06222021	MW15-09162020	MW-15-12172020
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	9/16/2020	12/17/2020	3/29/2021	6/22/2021	9/16/2020	12/17/2020
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds (cor	nt.)											
Phenanthrene	85-01-8	-					41				0.098 U	
Phenol	108-95-2	4,800					2.0 UJ				2.0 UJ	
Pyrene	129-00-0	240					1.7				0.20 U	
Total Naphthalenes		160					7,300				2.0 U	
Volatile Organic Compounds												
1,1,1,2-Tetrachloroethane	630-20-6											
1,1,1-Trichloroethane	71-55-6	200	1.0 U	1.0 U	4.0 U	4.0 U	1.0 U	1.0 U	4.0 U	4.0 U	1.0 U	1.0 U
1,1,2,2-Tetrachloroethane	79-34-5											
1,1,2-Trichloroethane	79-00-5											
1,1-Dichloroethane	75-34-3	7.7	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U
1,1-Dichloroethene	75-35-4	_										
1,1-Dichloropropene	563-58-6											
1,2,3-Trichlorobenzene	87-61-6	_					Ĭ					
1,2,3-Trichloropropane	96-18-4						İ					
1,2,4-Trichlorobenzene	120-82-1	_										
1,2,4-Trimethylbenzene	95-63-6	80	140	390	390	350	88	63	61	59	1.0 U	1.0 U
1,2-Dibromo-3-chloropropane	96-12-8									**		
1,2-Dibromoethane	106-93-4											
1,2-Dichlorobenzene	95-50-1											
1,2-Dichloroethane	107-06-2											
1,2-Dichloropropane	78-87-5											
1,3,5-Trimethylbenzene	108-67-8	80	40	120	130	110	24	18	13	13	1.0 U	1.0 U
1,3-Dichlorobenzene	541-73-1		10	120	100			10		13	1.00	1.0 0
1,3-Dichloropropane	142-28-9	_										
1,4-Dichlorobenzene	106-46-7											
2,2-Dichloropropane	594-20-7											
2,4-Dinitrotoluene	121-14-2											
2,6-Dinitrotoluene	606-20-2											
2-Chlorotoluene	95-49-8	_										
2-Hexanone	591-78-6											
2-Nitroaniline	88-74-4											
3-Nitroaniline	99-09-2											
4-Chloroaniline	106-47-8	-										
4-Chlorophenyl phenyl ether	7005-72-3	-										
4-Chlorotoluene	106-43-4	-										
4-Nitroaniline	100-01-6											
Acetone	67-64-1											
Benzene	71-43-2	5.0	8.8	9.3	6.1	11	1.0 U	1.0 U	4.4 U	4.4 U	1.0 U	1.0 U
Bis(2-chloroethyl)ether	111-44-4	J.0 	0.0	3.3	0.1		1.00	1.00	4.4 0	4.4 0	1.00	1.00
Bis(2-chloroisopropyl)ether	39638-32-9											
Bromobenzene	108-86-1						1		1			+
Bromochloromethane	74-97-5						1	1	1			1
Bromodichloromethane	75-27-4	-					1					
Bromoform	75-25-2						1					
Bromomethane	75-25-2						1		-			-
Carbon disulfide	74-83-9 75-15-0						1		-			-
Carbon disulide Carbon tetrachloride	75-13-0 56-23-5						 		-			-
	108-90-7						 					
Chlorobenzene							 					
Chloroethane Chloroform	75-00-3 67-66-3	70	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U



Table 5.1 Groundwater Monitoring Well Analytical Data 2010-2021 (µg/L)

		Location Name		MW-12	(cont.)			MW	-14		MV	V-15
		Sample Name	MW12-09152020	MW-12-12172020		MW12-06222021	MW14-09162020	MW-14-12172020	MW14-032921	MW14-06222021	MW15-09162020	MW-15-12172020
		Sample Date	9/15/2020	12/17/2020	3/29/2021	6/22/2021	9/16/2020	12/17/2020	3/29/2021	6/22/2021	9/16/2020	12/17/2020
Analyte	CAS No.	Screening Level (1)										
Volatile Organic Compounds (cont.)	-							L				
Chloromethane	74-87-3		2.0 U	2.0 U	7.5 U	7.5 U	2.0 U	2.0 U	7.5 U	7.5 U	2.0 U	2.0 U
cis-1,2-Dichloroethene	156-59-2	70	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U
cis-1,3-Dichloropropene	10061-01-5											
Cymene	99-87-6											
Dibromochloromethane	124-48-1											
Dibromomethane	74-95-3	-										
Dichlorodifluoromethane	75-71-8	-										
Ethylbenzene	100-41-4	700	410	1,100	1,100	1,100	40	42	42	54	1.0 U	1.0 U
Hexachlorobenzene	118-74-1	-										
Hexachloroethane	67-72-1											
Isopropylbenzene	98-82-8	800	54 J	1.0 U	55	51	12	9.8	7.5	9.8	1.0 U	1.0 U
Methyl ethyl ketone	78-93-3				15 U				15 U			
Methyl isobutyl ketone	108-10-1	640	1.3 U	1.3 U		13 U	1.3 U	1.3 U		13 U	1.3 U	1.3 U
Methylene chloride	75-09-2											
Methyl-tert-butyl ether	1634-04-4	-										
n-Butylbenzene	104-51-8											
Nitrobenzene	98-95-3	-										
n-Propylbenzene	103-65-1	800	20	27	20	17	6.1	4.6	5.0 U	5.0	1.0 U	1.0 U
sec-Butylbenzene	135-98-8											
Styrene	100-42-5	100	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U	5.0 U	5.0 U	1.0 U	1.0 U
tert-Butylbenzene	98-06-6											
Tetrachloroethene	127-18-4											
Toluene	108-88-3	1,000	130	360	400	380	3.0	3.2	7.5 U	7.5 U	1.0 U	1.0 U
trans-1,2-Dichloroethene	156-60-5											
trans-1,3-Dichloropropene	10061-02-6											
Trichloroethene	79-01-6	5.0	0.50 U	0.50 U	5.0 U	5.0 U	0.50 U	0.50 U	5.0 U	5.0 U	0.50 U	0.50 U
Trichlorofluoromethane	75-69-4											
Vinyl chloride	75-01-4		·									
Xylene (meta & para)	108-38-3/106-42-3		460	1,300	1,200	1,200	40	35	31	35	1.0 U	1.0 U
Xylene (ortho)	95-47-6		280	660	600	630	38	32	30	39	1.0 U	1.0 U
Xylene (total)	1330-20-7	1,000	740 ⁽⁵⁾	2,000	1,800 ⁽⁵⁾	1,800 ⁽⁵⁾	77 ⁽⁵⁾	68	61 ⁽⁵⁾	74 ⁽⁵⁾	1.0 U ⁽⁵⁾	1.0 U
Other												
PBDE-003	101-55-3		_									

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Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

-- Not available or not established.

BOLD Detected result exceeds criteria. The screening level is a proposed Site cleanup level.

- 16 Strongwards screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.

 2 A field duplicate was collected with this sample. The highest detected cressly as selected or compared to criteria. In feasible from the compared to the screening levels was prefered new as unjust by the proposed to the screening level for total naphthalenes.

 4 Naphthalene was analyzed by both USEPA 8260 only, the pay 6200 only the pay 6200 only t

Abbreviations:

CAS = Chemical Abstracts Service PBDE = Polybrominated diphenyl ethers

TEQ = Toxic equivalent cPAH = Carcinogenic polycyclic aromatic hydrocarbon

USEPA = U.S. Environmental Protection Agency

μg/L = Micrograms per liter MTCA = Model Toxics Control Act

Qualifiers:

J Analyte was detected and the concentration is estimated.

Modified from Floyd | Snider 2022 draft RI/FS.

- U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

	-	Location Name	MW-15	(cont.)			V-16			MW-	17	
		Sample Name	MW15-032921	MW15-06222021	MW16-09152020	MW-16-12172020	MW16-032921 ⁽²⁾	MW16-06222021	MW17-09152020	MW-17-12172020	MW17-032921	MW17-0622202
		Sample Date	3/29/2021	6/22/2021	9/15/2020	12/17/2020	3/29/2021	6/22/2021	9/15/2020	12/17/2020	3/29/2021	6/22/2021
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds												
1-Methylnaphthalene	90-12-0				1.3				0.10			
2,4,5-Trichlorophenol	95-95-4	1,600			2.0 U				2.0 U			
2,4,6-Trichlorophenol	88-06-2	8.0			2.0 U				2.0 U			
2,4-Dichlorophenol	120-83-2	48			2.0 U				2.0 U			
2,4-Dimethylphenol	105-67-9	320			0.98 U				1.0 U			
2,4-Dinitrophenol	51-28-5	32			2.0 UJ				2.0 UJ			
2-Chloronaphthalene	91-58-7	-										
2-Chlorophenol	95-57-8	40			0.98 U				1.0 U			
2-Methylnaphthalene	91-57-6				1.1				0.10 U			
2-Methylphenol	95-48-7	800			0.98 U				1.0 U			
2-Nitrophenol	88-75-5				2.0 U				2.0 U			
3- & 4-Methylphenol	15831-10-4	800			0.98 U				1.0 U			
4,6-Dinitro-o-cresol	534-52-1	1.3			49 U				5.0 UJ			
4-Chloro-3-methylphenol	59-50-7	1,600			4.9 U		1		5.0 U		1	
4-Nitrophenol	100-02-7	1,600			4.9 UJ				5.0 UJ			
Acenaphthene	83-32-9	480			0.80				0.20 U			
Acenaphthylene	208-96-8	_			0.20 U				0.20 U			
Acrylonitrile	107-13-1											
Anthracene	120-12-7	2,400			0.098 U				0.10 U			
Benzo(a)anthracene	56-55-3				0.20 U				0.20 U			
Benzo(a)pyrene	50-32-8				0.098 U				0.10 U			
Benzo(b)fluoranthene	205-99-2	-			0.20 U				0.20 U			
Benzo(g,h,i)perylene	191-24-2				0.20 U				0.20 U			
Benzo(k)fluoranthene	207-08-9				0.20 U				0.20 U			
Benzoic acid	65-85-0	-			0.20 0				0.20 0		1	
Benzyl alcohol	100-51-6										1	
Bis(2-chloroethoxy)methane	111-91-1										1	
Bis(2-ethylhexyl)phthalate	117-81-7	-									1	
Butyl benzyl phthalate	85-68-7											
Carbazole	86-74-8	-										
Chrysene	218-01-9				0.20 U				0.20 U			
Dibenzo(a,h)anthracene	53-70-3	-			0.20 U				0.20 U			
Indeno(1,2,3-c,d)pyrene	193-39-5				0.20 U				0.10 U			
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10			0.098 U		-		0.10 U		1	
cPAHs (MTCA TEQ-HallND)	BaPEq (U=1/2)	0.10			0.098 U		-		0.10 U		+	
Dibenzofuran	132-64-9	0.10			0.036 0		-		0.10 0		1	
Diethylphthalate	84-66-2						1		1		+	
Dimethyl phthalate	131-11-3						-				+	
	131-11-3 84-74-2						_				-	
Di-n-butyl phthalate	84-74-2 117-84-0						_				-	
Di-n-octyl phthalate	117-84-0 206-44-0	640			0.098 U		 		0.10 U		 	
Fluoranthene							 				 	
Fluorene	86-73-7	320			0.24		 		0.10 U		!	
Hexachlorobutadiene	87-68-3	-									-	
Hexachlorocyclopentadiene	77-47-4	-										
Isophorone	78-59-1		(2)	(2)		(2)	(2)	(2)	(2)	(2)	(2)	(2)
Naphthalene	91-20-3	160 ⁽³⁾	1.3 U ⁽³⁾	1.5 (3)	21	1,600 ⁽³⁾	3,100 ⁽³⁾	75 ⁽³⁾	2.0 U ⁽³⁾	70 ⁽³⁾	1.3 (3)	1.8 (3)
N-Nitroso-di-n-propylamine	621-64-7											
N-Nitrosodiphenylamine	86-30-6	-										
Pentachlorophenol	87-86-5	1.0			0.098 UJ				0.10 UJ			



Table 5.1 Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

Loc	cation Name			15 (cont.)			V-16			MW-		
		Sample Name	MW15-032921	MW15-06222021	MW16-09152020	MW-16-12172020	MW16-032921 (2)	MW16-06222021	MW17-09152020	MW-17-12172020	MW17-032921	MW17-06222021
		Sample Date	3/29/2021	6/22/2021	9/15/2020	12/17/2020	3/29/2021	6/22/2021	9/15/2020	12/17/2020	3/29/2021	6/22/2021
Analyte	CAS No.	Screening Level (1)										
Semivolatile Organic Compounds (con	nt.)											
Phenanthrene	85-01-8				0.098 U				0.10 U			
Phenol	108-95-2	4,800			2.0 UJ				2.0 UJ			
Pyrene	129-00-0	240			0.20 U				0.20 U			
Total Naphthalenes		160			23				0.10			
Volatile Organic Compounds												
1,1,1,2-Tetrachloroethane	630-20-6											
1,1,1-Trichloroethane	71-55-6	200	0.40 U	0.40 U	1.0 U	1.0 U	0.40 UJ	0.40 U	1.0 U	1.0 U	0.40 U	0.40 U
1,1,2,2-Tetrachloroethane	79-34-5											
1,1,2-Trichloroethane	79-00-5	-										
1,1-Dichloroethane	75-34-3	7.7	0.50 U	0.50 U	1.0 U	1.0 U	0.50 UJ	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U
1,1-Dichloroethene	75-35-4	-										
1,1-Dichloropropene	563-58-6											
1,2,3-Trichlorobenzene	87-61-6	-										
1,2,3-Trichloropropane	96-18-4											
1,2,4-Trichlorobenzene	120-82-1											
1,2,4-Trimethylbenzene	95-63-6	80	0.50 U	0.50 U	2.2	59	88	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U
1,2-Dibromo-3-chloropropane	96-12-8											
1,2-Dibromoethane	106-93-4											
1,2-Dichlorobenzene	95-50-1											
1,2-Dichloroethane	107-06-2											
1,2-Dichloropropane	78-87-5											
1,3,5-Trimethylbenzene	108-67-8	80	0.25 U	0.25 U	1.0 U	25	31	0.25 U	1.0 U	1.0 U	0.25 U	0.25 U
1,3-Dichlorobenzene	541-73-1											
1,3-Dichloropropane	142-28-9	_										
1,4-Dichlorobenzene	106-46-7											
2,2-Dichloropropane	594-20-7											
2,4-Dinitrotoluene	121-14-2	-										
2,6-Dinitrotoluene	606-20-2	_										
2-Chlorotoluene	95-49-8											
2-Hexanone	591-78-6	-										
2-Nitroaniline	88-74-4											
3-Nitroaniline	99-09-2											
4-Chloroaniline	106-47-8											
4-Chlorophenyl phenyl ether	7005-72-3	-										
4-Chlorotoluene	106-43-4											
4-Nitroaniline	100-01-6											
Acetone	67-64-1	-										
Benzene	71-43-2	5.0	0.44 U	0.44 U	1.0 U	2.7	1.9	0.44 U	1.0 U	1.0 U	0.44 U	0.44 U
Bis(2-chloroethyl)ether	111-44-4		2.110	2		2.7		50		1.0 0		
Bis(2-chloroisopropyl)ether	39638-32-9											
Bromobenzene	108-86-1	-										
Bromochloromethane	74-97-5											
Bromodichloromethane	75-27-4						 					
Bromoform	75-25-2											
Bromomethane	74-83-9				ł		 		1		1	
Carbon disulfide	75-15-0				1		 					
Carbon tetrachloride	56-23-5	-										
Chlorobenzene	108-90-7				 		1				1	
					-		-				1	
Chloroethane	75-00-3		0.50.11	0.5011	10.11	10	0.50.11	0.50.11	1011	10.1/	0.50.1/	0.5011
Chloroform	67-66-3	70	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U



Table 5.1
Groundwater Monitoring Well Analytical Data 2010–2021 (μg/L)

Analyte Volatile Organic Compounds (cont.) Chloromethane cis-1,2-Dichloroethene	CAS No. 74-87-3 156-59-2	Location Name Sample Name Sample Date Screening Level (1)	MW-15 MW15-032921 3/29/2021	MW15-06222021 6/22/2021	MW16-09152020 9/15/2020	MW-16-12172020	MW16-032921 ⁽²⁾	MW16-06222021	MW17-09152020	MW-1 MW-17-12172020	MW17-032921	MW17-06222021
Volatile Organic Compounds (cont.) Chloromethane cis-1,2-Dichloroethene	74-87-3	Sample Date Screening Level (1)					MW16-032921 (2)	MW16-06222021				
Volatile Organic Compounds (cont.) Chloromethane cis-1,2-Dichloroethene	74-87-3	Screening Level (1)	3/29/2021	6/22/2021	9/15/2020							
Volatile Organic Compounds (cont.) Chloromethane cis-1,2-Dichloroethene	74-87-3				3/13/2020	12/17/2020	3/29/2021	6/22/2021	9/15/2020	12/17/2020	3/29/2021	6/22/2021
Chloromethane cis-1,2-Dichloroethene												
cis-1,2-Dichloroethene												
	156-59-2		0.75 U	0.75 U	2.0 U	2.0 U	0.75 UJ	0.75 U	2.0 U	2.0 U	0.75 U	0.75 U
		70	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U
cis-1,3-Dichloropropene	10061-01-5											
Cymene	99-87-6											
Dibromochloromethane	124-48-1											
Dibromomethane	74-95-3											
Dichlorodifluoromethane	75-71-8											
Ethylbenzene	100-41-4	700	0.40 U	0.40 U	4.6	200	230	0.40 U	1.0 U	1.0 U	0.40 U	0.40 U
Hexachlorobenzene	118-74-1											
Hexachloroethane	67-72-1											
Isopropylbenzene	98-82-8	800	0.50 U	0.50 U	1.0 U	11	16	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U
Methyl ethyl ketone	78-93-3	-	1.5 U				1.5 U				1.5 U	
Methyl isobutyl ketone	108-10-1	640		1.3 U	1.3 U	1.3 U		1.3 U	1.3 U	1.3 U		1.3 U
Methylene chloride	75-09-2											
Methyl-tert-butyl ether	1634-04-4	-										
n-Butylbenzene	104-51-8	-										
Nitrobenzene	98-95-3	-										
n-Propylbenzene	103-65-1	800	0.50 U	0.50 U	1.0 U	3.3	5.0	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U
sec-Butylbenzene	135-98-8											
Styrene	100-42-5	100	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U	1.0 U	1.0 U	0.50 U	0.50 U
tert-Butylbenzene	98-06-6											
Tetrachloroethene	127-18-4											
Toluene	108-88-3	1,000	0.75 U	0.75 U	1.0 U	39	47	0.75 U	1.0 U	1.0 U	0.75 U	0.75 U
trans-1,2-Dichloroethene	156-60-5											
trans-1,3-Dichloropropene	10061-02-6											
Trichloroethene	79-01-6	5.0	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U	0.50 U
Trichlorofluoromethane	75-69-4											
Vinyl chloride	75-01-4											
Xylene (meta & para)	108-38-3/106-42-3		1.0 U	1.0 U	5.3	150	160	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Xylene (ortho)	95-47-6		0.50 U	0.50 U	3.3	120	180	0.86	1.0 U	1.0 U	0.50 U	0.50 U
Xylene (total)	1330-20-7	1,000	1.0 U ⁽⁵⁾	1.0 U ⁽⁴⁾	8.5 ⁽⁵⁾	260	340 ⁽⁵⁾	0.86	1.0 U ⁽⁵⁾	1.0 U	1.0 U ⁽⁵⁾	1.0 U ⁽⁵⁾
Other	= -	,	1.0 0				2.3		2.5 0			
PBDE-003	101-55-3											

Blank cells are intentional.

All criteria and results are rounded to two significant figures. Field parameters are not rounded.

- Not available or not established.

BOLD Detected result exceeds criteria. The screening level is a proposed Site cleanup level.

1 Groundwater screening levels for chemicals of interest are established in Table 5.2. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.

2 A field duplicate was collected with this sample. The highest detected result was selected and compared to criteria. If nether result was detected, the lowest RL was selected. 3 Where naphthalene was run by USEPA 8260 only, the naphthalene result is compared to the screening level for total naphthalenes.

4 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level. 5 Xylene (total) as calculated by Floyd | Snider was preferred over laboratory reported result.

Abbreviations:

CAS = Chemical Abstracts Service

PBDE = Polybrominated diphenyl ethers TEQ = Toxic equivalent

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

USEPA = U.S. Environmental Protection Agency

μg/L = Micrograms per liter MTCA = Model Toxics Control Act

Qualifiers:

J Analyte was detected and the concentration is estimated.

U Analyte is not detected at the associated reporting limit.

UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.2 Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-01	SB-02	SB-03	SB-04	SB-05	SB-	06	CP	-07	CE	3-08	S.D.	3-09
	Sample Name	SB1-34-GW	SB2-34-GW	SB3-33-GW	SB4-35-GW	SB5-35-GW	SB06-22-RGW	SB06-37-RGW	SB07-24-RGW	SB07-37-RGW	SB08-26-RGW	SB08-42-RGW	SB09-32-RGW	SB09-43-RGW
	Sample Date	1/29/2007	1/29/2007	1/31/2007	5/11/2007	5/11/2007	4/2/2009	4/2/2009	4/2/2009	4/2/2009	4/1/2009	4/2/2009	4/1/2009	4/1/2009
Analyte	CAS No.	1/23/2007	1/23/2007	1/31/2007	3/11/2007	3/11/2007	4/2/2003	4/2/2003	4/2/2003	4/2/2003	4/1/2003	4/2/2003	4/1/2003	4/1/2003
Semivolatile Organic Compounds	6/15/110/													
1-Methylnaphthalene	90-12-0	0.87	0.50	0.20	0.12 U	190								
2-Methylnaphthalene	91-57-6	0.95	0.37	0.26	0.12 U	310								
Acenaphthene	83-32-9	0.86	0.61	0.18	0.12 U	170								
Acenaphthylene	208-96-8	0.12 U	0.096 U	0.14 U	0.12 U	11 U								
Anthracene	120-12-7	0.13	0.11	0.17	0.12 U	30								
Benzo(a)anthracene	56-55-3	0.081	0.011	0.15	0.020	14								
Benzo(a)pyrene	50-32-8	0.095	0.014	0.12	0.14	6.3								
Benzo(b)fluoranthene	205-99-2	0.093	0.026	0.14	0.16	8.8								
Benzo(g,h,i)perylene	191-24-2	0.072	0.020	0.079	0.19	2.3								
Benzo(k)fluoranthene	207-08-9	0.032	0.0096 U	0.051	0.054	3.9								
Chrysene	218-01-9	0.10	0.020	0.17	0.11	13								
Dibenzo(a,h)anthracene	53-70-3	0.015	0.0096 U	0.021	0.039	1.1 U								
Indeno(1,2,3-c,d)pyrene	193-39-5	0.056	0.013	0.065	0.15	2.4								
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.12	0.020	0.16	0.18	9.4								
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.12	0.019	0.16	0.18	9.3								
Fluoranthene	206-44-0	0.16	0.096 U	0.39	0.12 U	71								
Fluorene	86-73-7	0.57	0.36	0.18	0.12 U	120								
Hexachlorobutadiene	87-68-3						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Naphthalene	91-20-3	5.7	1.0	0.76	0.12 U	1,200	1.0 U	1.3	1.0 U	1.0 U	1.0 U	1.0 U	11	4,500
Pentachlorophenol	87-86-5													
Phenanthrene	85-01-8	0.70	0.62	0.79	0.12 U	200								
Pyrene	129-00-0	0.21	0.096 U	0.42	0.12 U	51								
Total Naphthalenes		7.5	1.9	1.2	0.12 U	1,700								
Volatile Organic Compounds														
1,1,1,2-Tetrachloroethane	630-20-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,1,1-Trichloroethane	71-55-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,1,2,2-Tetrachloroethane	79-34-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,1,2-Trichloroethane	79-00-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,1-Dichloroethane	75-34-3						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,1-Dichloroethene	75-35-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,1-Dichloropropene	563-58-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,2,3-Trichlorobenzene	87-61-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,2,3-Trichloropropane	96-18-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,2,4-Trichlorobenzene	120-82-1						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,2,4-Trimethylbenzene	95-63-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	1.2	140
1,2-Dibromo-3-chloropropane	96-12-8						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
1,2-Dibromoethane	106-93-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,2-Dichlorobenzene	95-50-1						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,2-Dichloroethane	107-06-2						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,2-Dichloropropane	78-87-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,3,5-Trimethylbenzene	108-67-8						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.73	60
1,3-Dichlorobenzene	541-73-1						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
1,3-Dichloropropane	142-28-9						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U



Table 5.2
Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-01	SB-02	SB-03	SB-04	SB-05	SB-	06	SB	-07	SI	3-08	SB	3-09
	Sample Name	SB1-34-GW	SB2-34-GW	SB3-33-GW	SB4-35-GW	SB5-35-GW	SB06-22-RGW	SB06-37-RGW	SB07-24-RGW	SB07-37-RGW	SB08-26-RGW	SB08-42-RGW	SB09-32-RGW	SB09-43-RGW
	Sample Date	1/29/2007	1/29/2007	1/31/2007	5/11/2007	5/11/2007	4/2/2009	4/2/2009	4/2/2009	4/2/2009	4/1/2009	4/2/2009	4/1/2009	4/1/2009
Analyte	CAS No.				5, ==, ===.	0/ ==/ ==0	., _, _	.,_,_	., _, _	.,_,_	., _,	., _, _	., _, _	., _, _
Volatile Organic Compounds (cont.)														
1,4-Dichlorobenzene	106-46-7						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
2,2-Dichloropropane	594-20-7						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
2-Chloroethyl vinyl ether	110-75-8						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
2-Chlorotoluene	95-49-8						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
2-Hexanone	591-78-6						2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	200 U
4-Chlorotoluene	106-43-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Acetone	67-64-1						5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	500 U
Benzene	71-43-2						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Bromobenzene	108-86-1						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Bromochloromethane	74-97-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Bromodichloromethane	75-27-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Bromoform	75-25-2						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
Bromomethane	74-83-9						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Carbon disulfide	75-15-0						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Carbon tetrachloride	56-23-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Chlorobenzene	108-90-7						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Chloroethane	75-00-3						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
Chloroform	67-66-3						1.1	0.20 U	1.0	1.6	1.1	0.64	0.26	20 U
Chloromethane	74-87-3						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
cis-1,2-Dichloroethene	156-59-2						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
cis-1,3-Dichloropropene	10061-01-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Cymene	99-87-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.75	20 U
Dibromochloromethane	124-48-1						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Dibromomethane	74-95-3						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Dichlorodifluoromethane	75-71-8						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Ethylbenzene	100-41-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.25	20 U
Iodomethane	74-88-4						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
Isopropylbenzene	98-82-8						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.95	20 U
Methyl ethyl ketone	78-93-3						5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	5.0 U	500 U
Methyl isobutyl ketone	108-10-1						2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	200 U
Methylene chloride	75-09-2						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
Methyl-tert-butyl ether	1634-04-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
n-Butylbenzene	104-51-8						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	47
n-Propylbenzene	103-65-1						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.81	20 U
sec-Butylbenzene	135-98-8						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.66	20 U
Styrene	100-42-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
tert-Butylbenzene	98-06-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Tetrachloroethene	127-18-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Toluene	108-88-3						1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100 U
trans-1,2-Dichloroethene	156-60-5						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
trans-1,3-Dichloropropene	10061-02-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Trichloroethene	79-01-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Trichlorofluoromethane	75-69-4			l	I		0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U



Table 5.2 Reconnaissance Groundwater Analytical Data (µg/L)

	Location Name	SB-01	SB-02	SB-03	SB-04	SB-05	SB-	06	SB	-07	SE	3-08	SE	3-09
	Sample Name	SB1-34-GW	SB2-34-GW	SB3-33-GW	SB4-35-GW	SB5-35-GW	SB06-22-RGW	SB06-37-RGW	SB07-24-RGW	SB07-37-RGW	SB08-26-RGW	SB08-42-RGW	SB09-32-RGW	SB09-43-RGW
	Sample Date	1/29/2007	1/29/2007	1/31/2007	5/11/2007	5/11/2007	4/2/2009	4/2/2009	4/2/2009	4/2/2009	4/1/2009	4/2/2009	4/1/2009	4/1/2009
Analyte	CAS No.													
Volatile Organic Compounds (co	ont.)													
Vinyl acetate	108-05-4						2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	2.0 U	200 U
Vinyl chloride	75-01-4						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Xylene (meta & para)	108-38-3/106-42-3						0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	40 U
Xylene (ortho)	95-47-6						0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	20 U
Xylene (total)	1330-20-7	<u> </u>					0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	40 U

Note

Blank cells are intentional.

Analytical results are reported to two significant figures.

-- Not available or not established.

1 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

μg/L = Micrograms per liter

MTCA = Model Toxics Control Act

TEQ = Toxic equivalent

USEPA = U.S. Environmental Protection Agency

Qualifier



Table 5.2
Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-	10	SB-	-11	SB	-21	SB-22	SB-23	SB-24	SE	3-25
	Sample Name	SB10-29	SB10-44	SB11-29	SB11-44	WCD-SB21-24.0-120611	WCD-SB21-39.0-120611	WCD-SB22-39.0-120611	WCD-SB23-39.0-120611	WCD-SB24-39.0-120611	WCD-SB25-23.0-121211	WCD-SB25-37.0-121211
	Sample Date	5/12/2009	5/12/2009	5/12/2009	5/12/2009	12/6/2011	12/6/2011	12/6/2011	12/6/2011	12/6/2011	12/12/2011	12/12/2011
Analyte	CAS No.											
Semivolatile Organic Compounds												
1-Methylnaphthalene	90-12-0											
2-Methylnaphthalene	91-57-6											
Acenaphthene	83-32-9											
Acenaphthylene	208-96-8											
Anthracene	120-12-7											
Benzo(a)anthracene	56-55-3											
Benzo(a)pyrene	50-32-8											
Benzo(b)fluoranthene	205-99-2											
Benzo(g,h,i)perylene	191-24-2											
Benzo(k)fluoranthene	207-08-9											
Chrysene	218-01-9											
Dibenzo(a,h)anthracene	53-70-3											
Indeno(1,2,3-c,d)pyrene	193-39-5											
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)											
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)											
Fluoranthene	206-44-0											
Fluorene	86-73-7											
Hexachlorobutadiene	87-68-3	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Naphthalene	91-20-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	640	1.0 U	920	1.0 U	1.0 U	1.0 U
Pentachlorophenol	87-86-5											
Phenanthrene	85-01-8											
Pyrene	129-00-0											
Total Naphthalenes												
Volatile Organic Compounds												
1,1,1,2-Tetrachloroethane	630-20-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,1-Trichloroethane	71-55-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,2,2-Tetrachloroethane	79-34-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,2-Trichloroethane	79-00-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloroethane	75-34-3	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloroethene	75-35-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloropropene	563-58-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,3-Trichlorobenzene	87-61-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,3-Trichloropropane	96-18-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,4-Trichlorobenzene	120-82-1	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,4-Trimethylbenzene	95-63-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	19	1.0 U	25	1.0 U	1.0 U	1.0 U
1,2-Dibromo-3-chloropropane	96-12-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dibromoethane	106-93-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichlorobenzene	95-50-1	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichloroethane	107-06-2	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichloropropane	78-87-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,3,5-Trimethylbenzene	108-67-8	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	6.9	1.0 U	8.3	1.0 U	1.0 U	1.0 U
1,3-Dichlorobenzene	541-73-1	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,3-Dichloropropane	142-28-9	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U



Table 5.2 Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-	10	SB-	11	SB-	-21	SB-22	SB-23	SB-24	SE	3-25
	Sample Name	SB10-29	SB10-44	SB11-29	SB11-44	WCD-SB21-24.0-120611	WCD-SB21-39.0-120611	WCD-SB22-39.0-120611	WCD-SB23-39.0-120611	WCD-SB24-39.0-120611	WCD-SB25-23.0-121211	WCD-SB25-37.0-121211
	Sample Date	5/12/2009	5/12/2009	5/12/2009	5/12/2009	12/6/2011	12/6/2011	12/6/2011	12/6/2011	12/6/2011	12/12/2011	12/12/2011
Analyte	CAS No.	-,,	,,	-,, · ·	-,,	, 0,	, , ,		,,,	,,,		,,
Volatile Organic Compounds (cont.)			•									
1,4-Dichlorobenzene	106-46-7	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2,2-Dichloropropane	594-20-7	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2-Chloroethyl vinyl ether	110-75-8	1.0 U	1.0 U	1.0 U	1.0 U							
2-Chlorotoluene	95-49-8	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2-Hexanone	591-78-6	2.0 U	2.0 U	2.0 U	2.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
4-Chlorotoluene	106-43-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Acetone	67-64-1	5.0 U	5.0 U	5.0 U	5.0 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U
Benzene	71-43-2	0.20 U	0.72	0.20 U	0.20 U	1.0 U	14	1.0 U	3.1	1.0 U	1.0 U	1.0 U
Bromobenzene	108-86-1	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromochloromethane	74-97-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromodichloromethane	75-27-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromoform	75-25-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromomethane	74-83-9	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Carbon disulfide	75-15-0	0.20 U	0.41	0.20 U	0.20 U							
Carbon tetrachloride	56-23-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chlorobenzene	108-90-7	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloroethane	75-00-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloroform	67-66-3	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloromethane	74-87-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	2.0
cis-1,2-Dichloroethene	156-59-2	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
cis-1,3-Dichloropropene	10061-01-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Cymene	99-87-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dibromochloromethane	124-48-1	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dibromomethane	74-95-3	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dichlorodifluoromethane	75-71-8	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Ethylbenzene	100-41-4	0.20 U	1.3	0.20 U	0.20 U	1.0 U	91	1.0 U	85	1.0 U	1.0 U	1.0 U
Iodomethane	74-88-4	1.0 U	1.0 U	1.0 U	1.0 U							
Isopropylbenzene	98-82-8	0.20 U	0.20 U	0.20 U	0.25	1.0 U	3.2	1.0 U	6.4	1.0 U	1.0 U	1.0 U
Methyl ethyl ketone	78-93-3	5.0 U	5.0 U	5.0 U	5.0 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U
Methyl isobutyl ketone	108-10-1	2.0 U	2.0 U	2.0 U	2.0 U	1.0 U	6.3	1.0 U				
Methylene chloride	75-09-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Methyl-tert-butyl ether	1634-04-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
n-Butylbenzene	104-51-8	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
n-Propylbenzene	103-65-1	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.1	1.0 U	2.0	1.0 U	1.0 U	1.0 U
sec-Butylbenzene	135-98-8	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Styrene	100-42-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
tert-Butylbenzene	98-06-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Tetrachloroethene	127-18-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Toluene	108-88-3	1.0 U	1.3	1.8	1.5	1.4	39	1.0 U	13	1.0 U	1.0 U	1.0 U
trans-1,2-Dichloroethene	156-60-5	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
trans-1,3-Dichloropropene	10061-02-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Trichloroethene	79-01-6	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Trichlorofluoromethane	75-69-4	0.20 U	0.20 U	0.20 U	0.20 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U



Table 5.2 Reconnaissance Groundwater Analytical Data (µg/L)

	Location Name	SB-	-10	SB-	B-11 SB-21		SB-22	SB-23	SB-24	SB	3-25	
	Sample Name	SB10-29	SB10-44	SB11-29	SB11-44	WCD-SB21-24.0-120611	WCD-SB21-39.0-120611	WCD-SB22-39.0-120611	WCD-SB23-39.0-120611	WCD-SB24-39.0-120611	WCD-SB25-23.0-121211	WCD-SB25-37.0-121211
	Sample Date	5/12/2009	5/12/2009	5/12/2009	5/12/2009	12/6/2011	12/6/2011	12/6/2011	12/6/2011	12/6/2011	12/12/2011	12/12/2011
Analyte	CAS No.											
Volatile Organic Compounds (cont.)												
Vinyl acetate	108-05-4	2.0 U	2.0 U	2.0 U	2.0 U							
Vinyl chloride	75-01-4	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U	0.20 U
Xylene (meta & para)	108-38-3/106-42-3	0.40 U	0.40 U	0.40 U	0.40 U							
Xylene (ortho)	95-47-6	0.20 U	0.20 U	0.20 U	0.20 U							
Xylene (total)	1330-20-7	0.40 U	0.40 U	0.40 U	0.40 U	1.0 U	97	1.0 U	61	1.0 U	1.0 U	1.0 U

Blank cells are intentional.

Analytical results are reported to two significant figures.

-- Not available or not established.

1 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

μg/L = Micrograms per liter

MTCA = Model Toxics Control Act

TEQ = Toxic equivalent

USEPA = U.S. Environmental Protection Agency

Qualifier



 $Table \ 5.2$ Reconnaissance Groundwater Analytical Data (µg/L)

	j			1		1	1		1	
	Location Name	SB-		SB-		SB-28		3-29		3-31
	Sample Name	WCD-SB26-23.0-121211	WCD-SB26-37.0-121211	WCD-SB27-23.0-121211	WCD-SB27-37.0-121211	WCD-SB28-23.0-121211	WCD-SB29-25.0-120611	WCD-SB29-39.0-120611	WCD-SB31-22-26'	WCD-SB31-36-40'
	Sample Date	12/12/2011	12/12/2011	12/12/2011	12/12/2011	12/12/2011	12/6/2011	12/6/2011	11/12/2012	11/12/2012
Analyte	CAS No.									
Semivolatile Organic Compounds									•	T
1-Methylnaphthalene	90-12-0									
2-Methylnaphthalene	91-57-6									
Acenaphthene	83-32-9									
Acenaphthylene	208-96-8									
Anthracene	120-12-7									
Benzo(a)anthracene	56-55-3									
Benzo(a)pyrene	50-32-8									
Benzo(b)fluoranthene	205-99-2									
Benzo(g,h,i)perylene	191-24-2									
Benzo(k)fluoranthene	207-08-9									
Chrysene	218-01-9									
Dibenzo(a,h)anthracene	53-70-3									
Indeno(1,2,3-c,d)pyrene	193-39-5									
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)									
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)									
Fluoranthene	206-44-0									
Fluorene	86-73-7									
Hexachlorobutadiene	87-68-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Naphthalene	91-20-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	5,400	1,400	2,700
Pentachlorophenol	87-86-5									
Phenanthrene	85-01-8									
Pyrene	129-00-0									
Total Naphthalenes										
Volatile Organic Compounds										
1,1,1,2-Tetrachloroethane	630-20-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,1-Trichloroethane	71-55-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,2,2-Tetrachloroethane	79-34-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,2-Trichloroethane	79-00-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloroethane	75-34-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloroethene	75-35-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloropropene	563-58-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,3-Trichlorobenzene	87-61-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,3-Trichloropropane	96-18-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,4-Trichlorobenzene	120-82-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,4-Trimethylbenzene	95-63-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	140	9.0	60
1,2-Dibromo-3-chloropropane	96-12-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dibromoethane	106-93-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichlorobenzene	95-50-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichloroethane	107-06-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichloropropane	78-87-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,3,5-Trimethylbenzene	108-67-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	51	3.0	20
1,3-Dichlorobenzene	541-73-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,3-Dichloropropane	142-28-9	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U



Table 5.2
Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-	26	SB-	27	SB-28	SE	3-29	SE	-31
	Sample Name	WCD-SB26-23.0-121211	WCD-SB26-37.0-121211	WCD-SB27-23.0-121211	WCD-SB27-37.0-121211	WCD-SB28-23.0-121211	WCD-SB29-25.0-120611	WCD-SB29-39.0-120611	WCD-SB31-22-26'	WCD-SB31-36-40'
	Sample Date	12/12/2011	12/12/2011	12/12/2011	12/12/2011	12/12/2011	12/6/2011	12/6/2011	11/12/2012	11/12/2012
Analyte	CAS No.			• •						
Volatile Organic Compounds (cont.)									•	•
1,4-Dichlorobenzene	106-46-7	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2,2-Dichloropropane	594-20-7	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2-Chloroethyl vinyl ether	110-75-8									
2-Chlorotoluene	95-49-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2-Hexanone	591-78-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
4-Chlorotoluene	106-43-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Acetone	67-64-1	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U
Benzene	71-43-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	7.7	1.0 U	1.4
Bromobenzene	108-86-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromochloromethane	74-97-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromodichloromethane	75-27-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromoform	75-25-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromomethane	74-83-9	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Carbon disulfide	75-15-0									
Carbon tetrachloride	56-23-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chlorobenzene	108-90-7	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloroethane	75-00-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloroform	67-66-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloromethane	74-87-3	1.0 U	2.3	1.0 U	3.3	1.0	1.0 U	1.0 U	1.0 U	1.0 U
cis-1,2-Dichloroethene	156-59-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
cis-1,3-Dichloropropene	10061-01-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Cymene	99-87-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dibromochloromethane	124-48-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dibromomethane	74-95-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dichlorodifluoromethane	75-71-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Ethylbenzene	100-41-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	620	24	150
Iodomethane	74-88-4									
Isopropylbenzene	98-82-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	23	1.0	10
Methyl ethyl ketone	78-93-3	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U
Methyl isobutyl ketone	108-10-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Methylene chloride	75-09-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Methyl-tert-butyl ether	1634-04-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
n-Butylbenzene	104-51-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
n-Propylbenzene	103-65-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	10	1.0 U	4.0
sec-Butylbenzene	135-98-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Styrene	100-42-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
tert-Butylbenzene	98-06-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
<u>Tetrachloroethene</u>	127-18-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Toluene	108-88-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	95	3.5	20
trans-1,2-Dichloroethene	156-60-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
trans-1,3-Dichloropropene	10061-02-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Trichloroethene	79-01-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.2	1.0 U	1.0 U	1.0 U
Trichlorofluoromethane	75-69-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U



Table 5.2
Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB	SB-26		-27	SB-28	S	6B-29	SI	3-31
	Sample Name \	WCD-SB26-23.0-121211	WCD-SB26-37.0-121211	WCD-SB27-23.0-121211	WCD-SB27-37.0-121211	WCD-SB28-23.0-121211	WCD-SB29-25.0-120611	WCD-SB29-39.0-120611	WCD-SB31-22-26'	WCD-SB31-36-40'
	Sample Date	12/12/2011	12/12/2011	12/12/2011	12/12/2011	12/12/2011	12/6/2011	12/6/2011	11/12/2012	11/12/2012
Analyte	CAS No.									
Volatile Organic Compounds (con-	t.)									
Vinyl acetate	108-05-4									
Vinyl chloride	75-01-4	0.20 U 0 U								
Xylene (meta & para)	108-38-3/106-42-3									
Xylene (ortho)	95-47-6									
Xylene (total)	1330-20-7	1.0 U	740	42	340					

Blank cells are intentional.

Analytical results are reported to two significant figures.

-- Not available or not established.

1 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

μg/L = Micrograms per liter

MTCA = Model Toxics Control Act

TEQ = Toxic equivalent

USEPA = U.S. Environmental Protection Agency

Qualifier



 $Table \ 5.2$ Reconnaissance Groundwater Analytical Data (µg/L)

	Location Name	SE	3-32	S	B-33	SI	3-34	S	B-35	S	B-36		SB-37
	Sample Name	WCD-SB32-22-26'	WCD-SB32-36-40'	WCD-SB33-22-26'	WCD-SB33-36-40'	WCD-SB34-26-30'	WCD-SB34-39-43'	WCD-SB35-26-30'	WCD-SB35-39-43'	WCD-SB36-26-30'	WCD-SB36-39-43'	WCD-SB37-26-30	WCD-SB37-35-39'
	Sample Date		11/12/2012	11/12/2012	11/12/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012
Analyte	CAS No.												
Semivolatile Organic Compounds													
1-Methylnaphthalene	90-12-0												
2-Methylnaphthalene	91-57-6												
Acenaphthene	83-32-9												
Acenaphthylene	208-96-8												
Anthracene	120-12-7												
Benzo(a)anthracene	56-55-3												
Benzo(a)pyrene	50-32-8												
Benzo(b)fluoranthene	205-99-2												
Benzo(g,h,i)perylene	191-24-2												
Benzo(k)fluoranthene	207-08-9												
Chrysene	218-01-9												
Dibenzo(a,h)anthracene	53-70-3												
Indeno(1,2,3-c,d)pyrene	193-39-5												
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)												
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)												
Fluoranthene	206-44-0												
Fluorene	86-73-7												
Hexachlorobutadiene	87-68-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Naphthalene	91-20-3	1,100	1,600	1,300	2,300	280	540	320	1,800	880	97	1.0 U	19
Pentachlorophenol	87-86-5												
Phenanthrene	85-01-8												
Pyrene	129-00-0												
Total Naphthalenes													
Volatile Organic Compounds													
1,1,1,2-Tetrachloroethane	630-20-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,1-Trichloroethane	71-55-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,2,2-Tetrachloroethane	79-34-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,2-Trichloroethane	79-00-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloroethane	75-34-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloroethene	75-35-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloropropene	563-58-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,3-Trichlorobenzene	87-61-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,3-Trichloropropane	96-18-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,4-Trichlorobenzene	120-82-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2,4-Trimethylbenzene	95-63-6	25	100	10	45	7.9	61	5.4	67	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dibromo-3-chloropropane	96-12-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dibromoethane	106-93-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichlorobenzene	95-50-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichloroethane	107-06-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,2-Dichloropropane	78-87-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,3,5-Trimethylbenzene	108-67-8	9.0	40	2.0	11	2.0	18	1.8	23	1.0 U	1.0 U	1.0 U	1.0 U
1,3-Dichlorobenzene	541-73-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,3-Dichloropropane	142-28-9	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U



Table 5.2
Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-	32	S	3-33	SE	3-34	SB-	35	S	B-36	SI	3-37
	Sample Name	WCD-SB32-22-26'	WCD-SB32-36-40'	WCD-SB33-22-26'	WCD-SB33-36-40'	WCD-SB34-26-30'	WCD-SB34-39-43'	WCD-SB35-26-30'	WCD-SB35-39-43'	WCD-SB36-26-30'	WCD-SB36-39-43'	WCD-SB37-26-30	WCD-SB37-35-39'
	Sample Date	11/12/2012	11/12/2012	11/12/2012	11/12/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012
Analyte	CAS No.												
Volatile Organic Compounds (cont.)													
1,4-Dichlorobenzene	106-46-7	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2,2-Dichloropropane	594-20-7	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2-Chloroethyl vinyl ether	110-75-8												
2-Chlorotoluene	95-49-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
2-Hexanone	591-78-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
4-Chlorotoluene	106-43-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Acetone	67-64-1	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U
Benzene	71-43-2	1.0 U	2.0	1.0 U	1.0 U	1.0 U	2.7	1.3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromobenzene	108-86-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromochloromethane	74-97-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromodichloromethane	75-27-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromoform	75-25-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Bromomethane	74-83-9	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Carbon disulfide	75-15-0												
Carbon tetrachloride	56-23-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chlorobenzene	108-90-7	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloroethane	75-00-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloroform	67-66-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloromethane	74-87-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
cis-1,2-Dichloroethene	156-59-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
cis-1,3-Dichloropropene	10061-01-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Cymene	99-87-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dibromochloromethane	124-48-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dibromomethane	74-95-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Dichlorodifluoromethane	75-71-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Ethylbenzene	100-41-4	60	180	8.0	36	23	160	10	98	1.0 U	1.0 U	1.0 U	1.0 U
Iodomethane	74-88-4												
Isopropylbenzene	98-82-8	4.0	20	1.0	6.0	1.4	13	1.1	13	1.0 U	1.0 U	1.0 U	1.0 U
Methyl ethyl ketone	78-93-3	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U	10 U
Methyl isobutyl ketone	108-10-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Methylene chloride	75-09-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Methyl-tert-butyl ether	1634-04-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
n-Butylbenzene	104-51-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
n-Propylbenzene	103-65-1	1.0	7.0	1.0 U	3.0	1.0 U	3.5	1.0 U	5.3	1.0 U	1.0 U	1.0 U	1.0 U
sec-Butylbenzene	135-98-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Styrene	100-42-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
tert-Butylbenzene	98-06-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Tetrachloroethene	127-18-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Toluene	108-88-3	6.0	28	1.0 U	3.0	2.6	20	1.0	6.9	1.0 U	1.0 U	1.0 U	1.0 U
trans-1,2-Dichloroethene	156-60-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
trans-1,3-Dichloropropene	10061-02-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Trichloroethene	79-01-6	2.0	1.0 U	1.1	1.0 U	1.0 U	1.1						
Trichlorofluoromethane	75-69-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U

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Table 5.2 Reconnaissance Groundwater Analytical Data (µg/L)

	Location Name	SB-32		SB-33		SB-34		SB-	35	SB	3-36	SE	B-37
	Sample Name	WCD-SB32-22-26'	WCD-SB32-36-40'	WCD-SB33-22-26'	WCD-SB33-36-40'	WCD-SB34-26-30'	WCD-SB34-39-43'	WCD-SB35-26-30'	WCD-SB35-39-43'	WCD-SB36-26-30'	WCD-SB36-39-43'	WCD-SB37-26-30	WCD-SB37-35-39'
	Sample Date	11/12/2012	11/12/2012	11/12/2012	11/12/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012	11/19/2012
Analyte	CAS No.												
Volatile Organic Compounds (cont	.)												
Vinyl acetate	108-05-4												
Vinyl chloride	75-01-4	0.20 U 0.20 U											
Xylene (meta & para)	108-38-3/106-42-3												
Xylene (ortho)	95-47-6	•											
Xylene (total)	1330-20-7	110	400	9.0	47	36	310	15	180	1.0 U	1.0 U	1.0 U	1.0 U

Blank cells are intentional.

Analytical results are reported to two significant figures.

-- Not available or not established.

1 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

μg/L = Micrograms per liter MTCA Model Toxics Control Act

TEQ = Toxic equivalent

USEPA = U.S. Environmental Protection Agency

Qualifier:



Table 5.2
Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-	-38	SB	-39	MW-15	MW-16	MW-17
	Sample Name	WCD-SB38-26-30'	WCD-SB38-39-43'	WCD-SB39-26-30'	WCD-SB39-39-43'	MW-15-W-39-44FT	MW-16-W-39-44FT	MW-17-W-39-44FT
	Sample Date	11/19/2012	11/19/2012	11/19/2012	11/19/2012	1/23/2019	1/23/2019	1/23/2019
Analyte	CAS No.					_,,		_,,
Semivolatile Organic Compounds								
1-Methylnaphthalene	90-12-0					0.20	540	240
2-Methylnaphthalene	91-57-6					0.10 U	760	19
Acenaphthene	83-32-9					0.34	160	210
Acenaphthylene	208-96-8					0.10 U	19	6.8
Anthracene	120-12-7					0.10 U	5.9	5.8
Benzo(a)anthracene	56-55-3					0.10 U	0.10 U	0.099 U
Benzo(a)pyrene	50-32-8					0.10 U	0.10 U	0.099 U
Benzo(b)fluoranthene	205-99-2					0.10 U	0.10 U	0.099 U
Benzo(g,h,i)perylene	191-24-2					0.10 U	0.10 U	0.099 U
Benzo(k)fluoranthene	207-08-9					0.10 U	0.10 U	0.099 U
Chrysene	218-01-9					0.10 U	0.10 U	0.099 U
Dibenzo(a,h)anthracene	53-70-3					0.10 U	0.10 U	0.099 U
Indeno(1.2.3-c.d)pyrene	193-39-5					0.10 U	0.10 U	0.099 U
cPAHs (MTCA TEQ-HalfND)	BaPEg (U=1/2)					0.10 U	0.10 U	0.099 U
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)					0.10 U	0.10 U	0.099 U
Fluoranthene	206-44-0					0.10 U	2.3	1.5
Fluorene	86-73-7					0.10 U	67	28
Hexachlorobutadiene	87-68-3	1.0 U	1.0 U	1.0 U	1.0 U			
Naphthalene	91-20-3	1.0 U	7.8	1.0 U	1.0 U	1.6 ⁽¹⁾	13.000 ⁽¹⁾	2.900 ⁽¹⁾
Pentachlorophenol	87-86-5					0.10 U	0.10 U	0.099 U
Phenanthrene	85-01-8					0.10 U	52	66
Pyrene	129-00-0					0.10 U	1.4	0.69
Total Naphthalenes						1.8	14,000	3,100
Volatile Organic Compounds						1.0	11,000	3,100
1,1,1,2-Tetrachloroethane	630-20-6	1.0 U	1.0 U	1.0 U	1.0 U			
1.1.1-Trichloroethane	71-55-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1,2,2-Tetrachloroethane	79-34-5	1.0 U	1.0 U	1.0 U	1.0 U	210 0	2.0	2.0 0
1,1,2-Trichloroethane	79-00-5	1.0 U	1.0 U	1.0 U	1.0 U			
1,1-Dichloroethane	75-34-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
1,1-Dichloroethene	75-35-4	1.0 U	1.0 U	1.0 U	1.0 U	210 0	2.0	2.0 0
1.1-Dichloropropene	563-58-6	1.0 U	1.0 U	1.0 U	1.0 U			
1,2,3-Trichlorobenzene	87-61-6	1.0 U	1.0 U	1.0 U	1.0 U			
1.2.3-Trichloropropane	96-18-4	1.0 U	1.0 U	1.0 U	1.0 U			
1,2,4-Trichlorobenzene	120-82-1	1.0 U	1.0 U	1.0 U	1.0 U			
1.2.4-Trimethylbenzene	95-63-6	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	290	40
1,2-Dibromo-3-chloropropane	96-12-8	1.0 U	1.0 U	1.0 U	1.0 U	2.0 0		.,
1,2-Dibromoethane	106-93-4	1.0 U	1.0 U	1.0 U	1.0 U			
1.2-Dichlorobenzene	95-50-1	1.0 U	1.0 U	1.0 U	1.0 U			
1.2-Dichloroethane	107-06-2	1.0 U	1.0 U	1.0 U	1.0 U			
1,2-Dichloropropane	78-87-5	1.0 U	1.0 U	1.0 U	1.0 U			
1,3,5-Trimethylbenzene	108-67-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	100	1.0 U
1,3-Dichlorobenzene	541-73-1	1.0 U	1.0 U	1.0 U	1.0 U	2.0		2.0 0
1,3-Dichloropropane	142-28-9	1.0 U	1.0 U	1.0 U	1.0 U			



Table 5.2
Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-	38	SB	-39	MW-15	MW-16	MW-17
	Sample Name	WCD-SB38-26-30'	WCD-SB38-39-43'	WCD-SB39-26-30'	WCD-SB39-39-43'	MW-15-W-39-44FT	MW-16-W-39-44FT	MW-17-W-39-44FT
	Sample Date	11/19/2012	11/19/2012	11/19/2012	11/19/2012	1/23/2019	1/23/2019	1/23/2019
Analyte	CAS No.							
Volatile Organic Compounds (cont.)								
1,4-Dichlorobenzene	106-46-7	1.0 U	1.0 U	1.0 U	1.0 U			
2,2-Dichloropropane	594-20-7	1.0 U	1.0 U	1.0 U	1.0 U			
2-Chloroethyl vinyl ether	110-75-8							
2-Chlorotoluene	95-49-8	1.0 U	1.0 U	1.0 U	1.0 U			
2-Hexanone	591-78-6	1.0 U	1.0 U	1.0 U	1.0 U			
4-Chlorotoluene	106-43-4	1.0 U	1.0 U	1.0 U	1.0 U			
Acetone	67-64-1	10 U	10 U	10 U	10 U			
Benzene	71-43-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	21	2.8
Bromobenzene	108-86-1	1.0 U	1.0 U	1.0 U	1.0 U			
Bromochloromethane	74-97-5	1.0 U	1.0 U	1.0 U	1.0 U			
Bromodichloromethane	75-27-4	1.0 U	1.0 U	1.0 U	1.0 U			
Bromoform	75-25-2	1.0 U	1.0 U	1.0 U	1.0 U			
Bromomethane	74-83-9	1.0 U	1.0 U	1.0 U	1.0 U			
Carbon disulfide	75-15-0							
Carbon tetrachloride	56-23-5	1.0 U	1.0 U	1.0 U	1.0 U			
Chlorobenzene	108-90-7	1.0 U	1.0 U	1.0 U	1.0 U			
Chloroethane	75-00-3	1.0 U	1.0 U	1.0 U	1.0 U			
Chloroform	67-66-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
Chloromethane	74-87-3	1.0 U	1.0 U	1.0 U	1.0 U	2.0 U	2.0 U	2.0 U
cis-1,2-Dichloroethene	156-59-2	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
cis-1,3-Dichloropropene	10061-01-5	1.0 U	1.0 U	1.0 U	1.0 U			
Cymene	99-87-6	1.0 U	1.0 U	1.0 U	1.0 U			
Dibromochloromethane	124-48-1	1.0 U	1.0 U	1.0 U	1.0 U			
Dibromomethane	74-95-3	1.0 U	1.0 U	1.0 U	1.0 U			
Dichlorodifluoromethane	75-71-8	1.0 U	1.0 U	1.0 U	1.0 U			
Ethylbenzene	100-41-4	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1,100	100
Iodomethane	74-88-4							
Isopropylbenzene	98-82-8	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	41	18
Methyl ethyl ketone	78-93-3	10 U	10 U	10 U	10 U			
Methyl isobutyl ketone	108-10-1	1.0 U	1.0 U	1.0 U	1.0 U	1.3 U	1.3 U	1.3 U
Methylene chloride	75-09-2	1.0 U	1.0 U	1.0 U	1.0 U			
Methyl-tert-butyl ether	1634-04-4	1.0 U	1.0 U	1.0 U	1.0 U			
n-Butylbenzene	104-51-8	1.0 U	1.0 U	1.0 U	1.0 U			
n-Propylbenzene	103-65-1	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	20	1.0 U
sec-Butylbenzene	135-98-8	1.0 U	1.0 U	1.0 U	1.0 U			
Styrene	100-42-5	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U
tert-Butylbenzene	98-06-6	1.0 U	1.0 U	1.0 U	1.0 U			
Tetrachloroethene	127-18-4	1.0 U	1.0 U	1.0 U	1.0 U			
Toluene	108-88-3	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	330	20
trans-1,2-Dichloroethene	156-60-5	1.0 U	1.0 U	1.0 U	1.0 U			
trans-1,3-Dichloropropene	10061-02-6	1.0 U	1.0 U	1.0 U	1.0 U			
Trichloroethene	79-01-6	1.0 U	1.0 U	1.0 U	1.0 U	0.50 U	0.50 U	0.50 U
Trichlorofluoromethane	75-69-4	1.0 U	1.0 U	1.0 U	1.0 U			



Table 5.2 Reconnaissance Groundwater Analytical Data (μg/L)

	Location Name	SB-	38	SB	-39	MW-15	MW-16	MW-17
	Sample Name	WCD-SB38-26-30'	WCD-SB38-39-43'	WCD-SB39-26-30'	WCD-SB39-39-43'	MW-15-W-39-44FT	MW-16-W-39-44FT	MW-17-W-39-44FT
	Sample Date	11/19/2012	11/19/2012	11/19/2012	11/19/2012	1/23/2019	1/23/2019	1/23/2019
Analyte	CAS No.							
Volatile Organic Compounds (cont.)								
Vinyl acetate	108-05-4							
Vinyl chloride	75-01-4	0.20 U	0.20 U	0.20 U	0.20 U			
Xylene (meta & para)	108-38-3/106-42-3					1.0 U	1,100	54
Xylene (ortho)	95-47-6					1.0 U	600	68
Xylene (total)	1330-20-7	1.0 U	1,700	120				

Blank cells are intentional.

Analytical results are reported to two significant figures.

-- Not available or not established.

1 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

μg/L = Micrograms per liter MTCA Model Toxics Control Act

TEQ = Toxic equivalent

USEPA = U.S. Environmental Protection Agency

Qualifier



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name	В	-01	B-02	B-03	B-04	B-05	SP-01	SP-02	SP-03	SP-04	SP-05	SP-06
		Sample Name	B01-6.5	B01-16.5	B02-6.0	B03-5.5	B04-9.5	B05-9.0	SP01-7-8	SP02-1-2	SP03-3-4	SP04-3-4	SP05-3-4	SP06-3-4
		Sample Date	4/30/1992	4/30/1992	5/4/1992	5/4/1992	5/5/1992	5/5/1992	7/20/2006	7/20/2006	7/20/2006	7/20/2006	9/12/2006	9/12/2006
		Sample Depth (feet)	6.5-6.5	16.5-16.5	6.0-6.0	5.5-5.5	9.5-9.5	9.0-9.0	7.0-8.0	1.0-2.0	3.0-4.0	3.0-4.0	3.0-4.0	3.0-4.0
Analyte	CAS No.	Screening Level (1)												
Metals	•													
Antimony	7440-36-0	32												
Arsenic	7440-38-2	20												
Beryllium	7440-41-7	160												
Cadmium	7440-43-9	2.0												
Chromium	7440-47-3	2,000												
Copper	7440-50-8	3,200												
Lead	7439-92-1	250												
Mercury	7439-97-6	2.0												
Nickel	7440-02-0	1,600												
Selenium	7782-49-2	400												
Silver	7440-22-4	400												
Thallium	7440-28-0	0.80												
Zinc	7440-66-6	24,000							1			1	1	
Total Petroleum Hydrocarbons		= 1,000												
Diesel-range organics	DRO		25 U	25 U					I			I	T	
Gasoline-range organics	GRO		7.0	5.0 U										
Semivolatile Organic Compound			7.0	3.0 0										
1-Methylnaphthalene	90-12-0	1							Ι	I		I	1	
2,4,5-Trichlorophenol	95-95-4	8,000					0.013 U	0.011 U						
2,4,6-Trichlorophenol	88-06-2	80					0.0076 U	0.0066 U						
2,4-Dichlorophenol	120-83-2	240					0.038 U	0.033 U						
2,4-Dimethylphenol	105-67-9	1,600					0.030 0	0.033 0						
2,4-Dinitrophenol	51-28-5	160												
2-Chlorophenol	95-57-8	400					1.9 U	1.6 U						
2-Methylnaphthalene	91-57-6						1.5 0	1.0 0						
2-Methylphenol	95-48-7	4,000												
2-Nitrophenol	88-75-5													
4,6-Dinitro-o-cresol	534-52-1	6.4												1
4-Chloro-3-methylphenol	59-50-7	8,000												
	106-44-5													
4-Methylphenol 4-Nitrophenol	106-44-5	8,000												
·					0.20.11	0.20.11			9.9	2.5 U	2.5 U	21	0.0050.11	0.42
Acenaphthene	83-32-9	4,800			0.20 U 0.20 U	0.20 U 0.20 U			3.9		2.5 0	31 5.0 U	0.0050 U 0.0055	0.42
Acenaphthylene	208-96-8	24.000				0.20 0				2.5 U				3.3
Anthracene	120-12-7	24,000			0.29				28	2.5 U	170	150	0.0076	5.5
Benzo(a)anthracene	56-55-3				0.88	0.51			33	15	52	55	0.023	6.5
Benzo(a)pyrene	50-32-8	0.10			1.5	0.81			45	24	93	50	0.038	5.9
Benzo(b)fluoranthene	205-99-2				0.71	0.41			67	47	87	62	0.066	5.4
Benzo(k)fluoranthene	207-08-9				0.37	0.22			18	15	34	24	0.019	1.6
Chrysene	218-01-9				1	0.6			45	25	240	99	0.037	6.8
Dibenzo(a,h)anthracene	53-70-3				0.68	0.040 U			7.0	5.8	18	6.9	0.0061	0.89
Indeno(1,2,3-c,d)pyrene	193-39-5				0.81	0.41			25	17	49	20	0.019	2.5
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10			1.9	0.97			60	34	120	68	0.052	7.7
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10			1.9	0.97			60	34	120	68	0.052	7.7
Benzo(g,h,i)perylene	191-24-2				0.96	0.46			28	21	50	20	0.020	2.9
Fluoranthene	206-44-0	3,200			3	1.9			76	11	50	180	0.028	11



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name	В	-01	B-02	B-03	B-04	B-05	SP-01	SP-02	SP-03	SP-04	SP-05	SP-06
		Sample Name	B01-6.5	B01-16.5	B02-6.0	B03-5.5	B04-9.5	B05-9.0	SP01-7-8	SP02-1-2	SP03-3-4	SP04-3-4	SP05-3-4	SP06-3-4
		Sample Date	4/30/1992	4/30/1992	5/4/1992	5/4/1992	5/5/1992	5/5/1992	7/20/2006	7/20/2006	7/20/2006	7/20/2006	9/12/2006	9/12/2006
		Sample Depth (feet)	6.5–6.5	16.5-16.5	6.0-6.0	5.5-5.5	9.5-9.5	9.0-9.0	7.0-8.0	1.0-2.0	3.0-4.0	3.0-4.0	3.0-4.0	3.0-4.0
Analyte	CAS No.	Screening Level (1)												
Semivolatile Organic Compound	s (cont.)													
Fluorene	86-73-7	3,200			0.20	0.12			4.7	2.5 U	3.4	29	0.0050 U	3.5
Naphthalene	91-20-3	5.0 ⁽²⁾			0.27 (2)	0.098 U (2)			5.9 ⁽²⁾	29 ⁽²⁾	2.6 (2)	5.0 U ⁽²⁾	0.0050 U (2)	0.25 U ⁽²⁾
Pentachlorophenol	87-86-5	2.5					0.0025 U	0.0022 U	25 UJ	25 UJ	40 UJ	70 UJ		
Phenanthrene	85-01-8				2.1	1.4			27	3.6	9.8	86	0.013	16
Phenol	108-95-2	24,000												
Pyrene	129-00-0	2,400			3.6	0.13			87	14	60	150	0.089	16
Tetrachlorophenols (total)	25167-83-3						0.0076 U	0.0066 U						
Total Naphthalenes		5.0												
Volatile Organic Compounds									•					
1,1,1-Trichloroethane	71-55-6	2												
1,1-Dichloroethane	75-34-3	180												
1,2,4-Trimethylbenzene	95-63-6	1.3												
1,2-Dichlorobenzene	95-50-1		0.029 U	0.026 U										
1,3,5-Trimethylbenzene	108-67-8	1.3												
1,3-Dichlorobenzene	541-73-1		0.029 U	0.026 U										
1,4-Dichlorobenzene	106-46-7		0.029 U	0.026 U										
Benzene	71-43-2	0.030	0.029 U	0.026 U										
Chlorobenzene	108-90-7		0.029 U	0.026 U										
Chloroform	67-66-3	32												
Chloromethane	74-87-3													
cis-1,2-Dichloroethene	156-59-2	160												
Ethylbenzene	100-41-4	6.0	0.029 U	0.026 U										
Isopropylbenzene	98-82-8	8,000												
Methyl isobutyl ketone	108-10-1	6,400												
n-Propylbenzene	103-65-1	8,000												
Styrene	100-42-5	16,000												
Toluene	108-88-3	7.0	0.029 U	0.026 U										
Trichloroethene	79-01-6	0.030												
Xylene (meta & para)	108-38-3/106-42-3													
Xylene (ortho)	95-47-6													
Xylene (total)	1330-20-7	9.0	0.029 U	0.026 U										

Blank cells are intentional.

All criteria and results are rounded to two significant figures.

-- Not available or not established.

BOLD Detected result exceeds criteria. The screening level is a proposed Site cleanup level.

- 1 Soil screening levels for all compounds designated as chemicals of interest are presented in Table 5.5. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.
- 2 Where naphthalene was run by USEPA 8260 only, or methylnaphthalenes were not run by USEPA 8270, the naphthalene result is compared to the screening level for total naphthalenes.
- 3 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level.

Abbreviation

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

mg/kg = Milligrams per kilogram

MTCA = Model Toxics Control Act TEQ Toxic equivalent

USEPA = U.S. Environmental Protection Agency

- J Analyte was detected and the concentration is estimated.
- JB Concentration is estimated due to presence of blank contamination.
- U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name	SP-07	SP-08	SP-09	SS-01	SS-02	SS-03	SS-04	SS-05	SS-06	MW			W-08
		Sample Name	SP07-7-8	SP08-3-4	SP09-7-8	SS-01	SS-02	SS-03	SS-04	SS-05	SS-06	MW7-17.5-19	MW7-25-26.5	MW8-15-16.5	MW8-25-26.5
		Sample Date	9/12/2006	9/12/2006	9/12/2006	9/21/2006	9/21/2006	9/21/2006	9/21/2006	9/21/2006	9/21/2006	1/26/2007	1/26/2007	1/31/2007	1/31/2007
Analista	CACNIC	Sample Depth (feet)	7.0–8.0	3.0-4.0	7.0-8.0	0.5–1.0	0.5-1.0	0.5–1.0	0.5-1.0	0.5-1.0	0.5-1.0	17.5–19.0	25.0-26.5	15.0–16.5	25.0-26.5
Analyte	CAS No.	Screening Level (1)					<u> </u>		<u> </u>				<u> </u>		
Metals	7440.25.0	22		1	1		1		T	1	T	1	<u> </u>		I
Antimony	7440-36-0	32													
Arsenic	7440-38-2	20													
Beryllium	7440-41-7	160													
Cadmium	7440-43-9	2.0					070								
Chromium	7440-47-3	2,000				9.2	370	9.3	8.5	8.0	18				
Copper	7440-50-8	3,200								0.0	100				
Lead	7439-92-1	250				68	22	11	11	26	190				
Mercury	7439-97-6	2.0													
Nickel	7440-02-0	1,600													
Selenium	7782-49-2	400											ļ		
Silver	7440-22-4	400													
Thallium	7440-28-0	0.80													
Zinc	7440-66-6	24,000		<u> </u>			<u> </u>		<u> </u>	<u> </u>	<u> </u>	<u> </u>	<u> </u>		<u> </u>
Total Petroleum Hydrocarbons	_			•	•					•		•			1
Diesel	DRO														
Gasoline	GRO														
Semivolatile Organic Compounds															
1-Methylnaphthalene	90-12-0											1.5	70	0.0069 U	0.0069 U
2,4,5-Trichlorophenol	95-95-4	8,000													
2,4,6-Trichlorophenol	88-06-2	80													
2,4-Dichlorophenol	120-83-2	240													
2,4-Dimethylphenol	105-67-9	1,600													
2,4-Dinitrophenol	51-28-5	160													
2-Chlorophenol	95-57-8	400													
2-Methylnaphthalene	91-57-6											2.8	130	0.0069 U	0.0069 U
2-Methylphenol	95-48-7	4,000													
2-Nitrophenol	88-75-5														
4,6-Dinitro-o-cresol	534-52-1	6.4													
4-Chloro-3-methylphenol	59-50-7	8,000													
4-Methylphenol	106-44-5	8,000													
4-Nitrophenol	100-02-7														
Acenaphthene	83-32-9	4,800	0.25 U	0.25 U	0.0050 U							3.9	49	0.0085	0.0069 U
Acenaphthylene	208-96-8		0.26	0.52	0.0050 U							0.12	2.7	0.061	0.0069 U
Anthracene	120-12-7	24,000	0.52	1.0	0.0050 U							1.5	16	0.098	0.0069 U
Benzo(a)anthracene	56-55-3		2.1	1.4	0.0050 U							1.4	6.7	0.33	0.0069 U
Benzo(a)pyrene	50-32-8	0.10	6.4	2.3	0.0050 U							0.66	3.8	0.29	0.0069 U
Benzo(b)fluoranthene	205-99-2		9.6	3.0	0.0050 U							0.92	4.9	0.83	0.011
Benzo(k)fluoranthene	207-08-9		2.3	0.87	0.0050 U							0.34	2.0	0.26	0.0069 U
Chrysene	218-01-9		5.2	2.4	0.0050 U							1.1	7.1	0.47	0.0069 U
Dibenzo(a,h)anthracene	53-70-3		1.1	0.67	0.0050 U							0.092	0.54	0.13	0.0069 U
Indeno(1,2,3-c,d)pyrene	193-39-5		2.9	2.4	0.0050 U							0.25	1.4	0.37	0.0069 U
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	8.3	3.2	0.0050 U		1		1	1	1	0.97	5.4	0.49	0.0060
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	8.3	3.2	0.0050 U							0.97	5.4	0.49	0.0011
Benzo(g,h,i)perylene	191-24-2		3.1	3.1	0.0050 U							0.27	1.5	0.39	0.0069 U
Fluoranthene	206-44-0	3,200	3.3	2.1	0.0050 U		1		1	 	 	6.1	35	0.42	0.0069 U



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name	SP-07	SP-08	SP-09	SS-01	SS-02	SS-03	SS-04	SS-05	SS-06	MW	-07	M	W-08
		Sample Name	SP07-7-8	SP08-3-4	SP09-7-8	SS-01	SS-02	SS-03	SS-04	SS-05	SS-06	MW7-17.5-19	MW7-25-26.5	MW8-15-16.5	MW8-25-26.5
		Sample Date	9/12/2006	9/12/2006	9/12/2006	9/21/2006	9/21/2006	9/21/2006	9/21/2006	9/21/2006	9/21/2006	1/26/2007	1/26/2007	1/31/2007	1/31/2007
		Sample Depth (feet)	7.0-8.0	3.0-4.0	7.0-8.0	0.5-1.0	0.5-1.0	0.5-1.0	0.5-1.0	0.5-1.0	0.5-1.0	17.5-19.0	25.0-26.5	15.0-16.5	25.0-26.5
Analyte	CAS No.	Screening Level (1)													
Semivolatile Organic Compound	s (cont.)														
Fluorene	86-73-7	3,200	0.25 U	0.25 U	0.0050 U							4.0	39	0.012	0.0069 U
Naphthalene	91-20-3	5.0 ⁽²⁾	0.25 U ⁽²⁾	0.25 U ⁽²⁾	0.0050 U ⁽²⁾							2.0	150	0.0081	0.0069 U
Pentachlorophenol	87-86-5	2.5													
Phenanthrene	85-01-8		0.25 U	1.5	0.0050 U							11	90	0.15	0.0069 U
Phenol	108-95-2	24,000													
Pyrene	129-00-0	2,400	5.9	3.0	0.0050 U							4.8	32	0.42	0.0069 U
Tetrachlorophenols (total)	25167-83-3														
Total Naphthalenes		5.0										6.3	350	0.0081	0.0069 U
Volatile Organic Compounds															
1,1,1-Trichloroethane	71-55-6	2													
1,1-Dichloroethane	75-34-3	180													
1,2,4-Trimethylbenzene	95-63-6	1.3													
1,2-Dichlorobenzene	95-50-1														
1,3,5-Trimethylbenzene	108-67-8	1.3													
1,3-Dichlorobenzene	541-73-1														
1,4-Dichlorobenzene	106-46-7														
Benzene	71-43-2	0.030													
Chlorobenzene	108-90-7														
Chloroform	67-66-3	32													
Chloromethane	74-87-3														
cis-1,2-Dichloroethene	156-59-2	160													ĺ
Ethylbenzene	100-41-4	6.0													ĺ
Isopropylbenzene	98-82-8	8,000													ĺ
Methyl isobutyl ketone	108-10-1	6,400													
n-Propylbenzene	103-65-1	8,000													
Styrene	100-42-5	16,000													
Toluene	108-88-3	7.0													
Trichloroethene	79-01-6	0.030				_									
Xylene (meta & para)	108-38-3/106-42-3														
Xylene (ortho)	95-47-6														
Xylene (total)	1330-20-7	9.0													

Blank cells are intentional.

All criteria and results are rounded to two significant figures.

-- Not available or not established. **BOLD** Detected result exceeds criteria.

The screening level is a proposed Site cleanup level.

- 1 Soil screening levels for all compounds designated as chemicals of interest are presented in Table 5.5. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.
- 2 Where naphthalene was run by USEPA 8260 only, or methylnaphthalenes were not run by USEPA 8270, the naphthalene result is compared to the screening level for total naphthalenes.
- 3 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

mg/kg = Milligrams per kilogram

MTCA = Model Toxics Control Act TEQ Toxic equivalent

USEPA = U.S. Environmental Protection Agency

Qualifiers:

J Analyte was detected and the concentration is estimated.

JB Concentration is estimated due to presence of blank contamination.

 $\ensuremath{\mathsf{U}}$ Analyte is not detected at the associated reporting limit.

UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name		MV	V-09		s	B-01	SI	3-02	SE	3-03	SB	-04	SB	-05
		Sample Name	MW9-24-25	MW9-45-47	MW9-55.5-57.5	MW9-68-70	SB1-9-10	SB1-19-20.5	SB2-14-15.5	SB2-19-20.5	SB3-10-11.5	SB3-15-16.5	SB4-10-11.5	SB4-20-21.5	SB5-15-16.5	SB5-35-35-36.5
		Sample Date	9/6/2007	9/6/2007	9/6/2007	9/6/2007	1/29/2007	1/29/2007	1/29/2007	1/29/2007	1/31/2007	1/31/2007	5/11/2007	5/11/2007	5/11/2007	5/11/2007
		Sample Depth (feet)	24.0-25.0	45.0-47.0	55.5-57.5	68.0-70.0	9.0-10.0	19.0-20.5	14.0-15.5	19.0-20.5	10.0-11.5	15.0-16.5	10.0-11.5	20.0-21.5	15.0-16.5	35.0-36.5
Analyte	CAS No.	Screening Level (1)														
Metals																
Antimony	7440-36-0	32														
Arsenic	7440-38-2	20														
Beryllium	7440-41-7	160														
Cadmium	7440-43-9	2.0														
Chromium	7440-47-3	2,000														
Copper	7440-50-8	3,200														
Lead	7439-92-1	250														
Mercury	7439-97-6	2.0														
Nickel	7440-02-0	1,600														
Selenium	7782-49-2	400														
Silver	7440-22-4	400														
Thallium	7440-28-0	0.80														
Zinc	7440-66-6	24,000														
Total Petroleum Hydrocarbons																
Diesel	DRO															
Gasoline	GRO															
Semivolatile Organic Compounds																
1-Methylnaphthalene	90-12-0		0.55	0.16	0.0072 U	0.0085 U	0.0072 U	0.0072 U	0.0074 U	0.0077 U	0.47	0.066	0.0081 U	0.0077 U	0.0072 U	0.19
2,4,5-Trichlorophenol	95-95-4	8,000														
2,4,6-Trichlorophenol	88-06-2	80														
2,4-Dichlorophenol	120-83-2	240														
2,4-Dimethylphenol	105-67-9	1,600														
2,4-Dinitrophenol	51-28-5	160														
2-Chlorophenol	95-57-8	400														
2-Methylnaphthalene	91-57-6		0.031	0.16	0.0072 U	0.0085 U	0.0072 U	0.0072 U	0.0074 U	0.0077 U	0.54	0.070	0.0081 U	0.0077 U	0.0072 U	0.16
2-Methylphenol	95-48-7	4,000														
2-Nitrophenol	88-75-5															
4,6-Dinitro-o-cresol	534-52-1	6.4														
4-Chloro-3-methylphenol	59-50-7	8,000														
4-Methylphenol	106-44-5	8,000														
4-Nitrophenol	100-02-7															
Acenaphthene	83-32-9	4,800	3.9	0.029	0.0072 U	0.0085 U	0.0072 U	0.011	0.0074 U	0.0077 U	0.75	0.078	0.0081 U	0.0077 U	0.0072 U	0.23
Acenaphthylene	208-96-8		0.093	0.0093 U	0.0072 U	0.0085 U	0.0099	0.0084	0.011	0.0077 U	0.52	0.10	0.011	0.0077 U	0.0072 U	0.0084
Anthracene	120-12-7	24,000	7.0	0.0093 U	0.0072 U	0.0085 U	0.0087	0.027	0.0074 U	0.0077 U	3.3	0.30	0.015	0.0077 U	0.0072 U	0.066
Benzo(a)anthracene	56-55-3		2.1	0.0093 U	0.0072 U	0.0085 U	0.025	0.050	0.0074 U	0.0077 U	5.3	0.49	0.012	0.0077 U	0.0072 U	0.0083 U
Benzo(a)pyrene	50-32-8	0.10	0.83	0.0093 U	0.0072 U	0.0085 U	0.030	0.053	0.0074 U	0.0077 U	5.4	0.49	0.11	0.017	0.0072 U	0.0083 U
Benzo(b)fluoranthene	205-99-2		1.1	0.0093 U	0.0072 U	0.0085 U	0.030	0.051	0.0074 U	0.0077 U	6.3	0.55	0.093	0.016	0.0072 U	0.0083 U
Benzo(k)fluoranthene	207-08-9		0.46	0.0093 U	0.0072 U	0.0085 U	0.0099	0.017	0.0074 U	0.0077 U	2.1	0.22	0.054	0.0099	0.0072 U	0.0083 U
Chrysene	218-01-9		2.5	0.0093 U	0.0072 U	0.0085 U	0.030	0.060	0.0074 U	0.0077 U	6.5	0.59	0.068	0.010	0.0072 U	0.0083 U
Dibenzo(a,h)anthracene	53-70-3		0.15	0.0093 U	0.0072 U	0.0085 U	0.0072 U	0.0082	0.0074 U	0.0077 U	0.85	0.086	0.036	0.0077 U	0.0072 U	0.0083 U
Indeno(1,2,3-c,d)pyrene	193-39-5		0.33	0.0093 U	0.0072 U	0.0085 U	0.016	0.024	0.0074 U	0.0077 U	2.8	0.26	0.12	0.017	0.0072 U	0.0083 U
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	1.3	0.0093 U	0.0072 U	0.0085 U	0.039	0.069	0.0074 U	0.0077 U	7.2	0.66	0.14	0.022	0.0072 U	0.0083 U
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	1.3	0.0093 U	0.0072 U	0.0085 U	0.038	0.069	0.0074 U	0.0077 U	7.2	0.66	0.14	0.021	0.0072 U	0.0083 U
Benzo(g,h,i)perylene	191-24-2		0.32	0.0093 U	0.0072 U	0.0085 U	0.021	0.031	0.0074 U	0.0077 U	3.4	0.30	0.15	0.021	0.0072 U	0.0083 U
Fluoranthene	206-44-0	3,200	9.5	0.0093 U	0.011	0.0085 U	0.038	0.084	0.0074 U	0.0077 U	10	0.98	0.0081 U	0.0077 U	0.0072 U	0.12



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name		MV	/-09		SI	B-01	SI	3-02	SE	3-03	SB	-04	SB	3-05
		Sample Name	MW9-24-25	MW9-45-47	MW9-55.5-57.5	MW9-68-70	SB1-9-10	SB1-19-20.5	SB2-14-15.5	SB2-19-20.5	SB3-10-11.5	SB3-15-16.5	SB4-10-11.5	SB4-20-21.5	SB5-15-16.5	SB5-35-35-36.5
		Sample Date	9/6/2007	9/6/2007	9/6/2007	9/6/2007	1/29/2007	1/29/2007	1/29/2007	1/29/2007	1/31/2007	1/31/2007	5/11/2007	5/11/2007	5/11/2007	5/11/2007
		Sample Depth (feet)	24.0-25.0	45.0-47.0	55.5-57.5	68.0-70.0	9.0-10.0	19.0-20.5	14.0-15.5	19.0-20.5	10.0-11.5	15.0-16.5	10.0-11.5	20.0-21.5	15.0-16.5	35.0-36.5
Analyte	CAS No.	Screening Level (1)														
Semivolatile Organic Compounds (cont.)															
Fluorene	86-73-7	3,200	2.9	0.0093 U	0.0072 U	0.0085 U	0.0072 U	0.0085	0.0074 U	0.0077 U	1.2	0.16	0.0081 U	0.0077 U	0.0072 U	0.19
Naphthalene	91-20-3	5.0 ⁽²⁾	0.064	3.8	0.0072 U	0.0085 U	0.0072 U	0.0072 U	0.0074 U	0.0077 U	0.55	0.041	0.0081 U	0.0077 U	0.0072 U	1.4
Pentachlorophenol	87-86-5	2.5														
Phenanthrene	85-01-8		23	0.0093 U	0.025	0.0085 U	0.024	0.097	0.0074 U	0.0077 U	11	1.0	0.0081 U	0.0077 U	0.0072 U	0.54
Phenol	108-95-2	24,000														
Pyrene	129-00-0	2,400	7.6	0.0093 U	0.0097	0.0085 U	0.059	0.15	0.0074 U	0.0077 U	15	1.3	0.0099	0.0077 U	0.0072 U	0.080
Tetrachlorophenols (total)	25167-83-3															
Total Naphthalenes		5.0	0.64	4.1	0.0072 U	0.0085 U	0.0072 U	0.0072 U	0.0074 U	0.0077 U	1.6	0.18	0.0081 U	0.0077 U	0.0072 U	1.8
Volatile Organic Compounds																
1,1,1-Trichloroethane	71-55-6	2														
1,1-Dichloroethane	75-34-3	180														
1,2,4-Trimethylbenzene	95-63-6	1.3														
1,2-Dichlorobenzene	95-50-1															
1,3,5-Trimethylbenzene	108-67-8	1.3														
1,3-Dichlorobenzene	541-73-1															
1,4-Dichlorobenzene	106-46-7															
Benzene	71-43-2	0.030														
Chlorobenzene	108-90-7															
Chloroform	67-66-3	32														
Chloromethane	74-87-3															
cis-1,2-Dichloroethene	156-59-2	160														
Ethylbenzene	100-41-4	6.0														
Isopropylbenzene	98-82-8	8,000														
Methyl isobutyl ketone	108-10-1	6,400														
n-Propylbenzene	103-65-1	8,000														
Styrene	100-42-5	16,000														
Toluene	108-88-3	7.0														
Trichloroethene	79-01-6	0.030														
Xylene (meta & para)	108-38-3/106-42-3															
Xylene (ortho)	95-47-6															
Xylene (total)	1330-20-7	9.0														

Blank cells are intentional.

All criteria and results are rounded to two significant figures.

-- Not available or not established.

BOLD Detected result exceeds criteria. The screening level is a proposed Site cleanup level.

- 1 Soil screening levels for all compounds designated as chemicals of interest are presented in Table 5.5. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.
- 2 Where naphthalene was run by USEPA 8260 only, or methylnaphthalenes were not run by USEPA 8270, the naphthalene result is compared to the screening level for total naphthalenes.
- 3 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level.

Abbreviatio

CAS = Chemical Abstracts Service cPAH = Carcinogenic polycyclic aromatic hydrocarbon mg/kg = Milligrams per kilogram MTCA = Model Toxics Control Act TEQ Toxic equivalent

USEPA = U.S. Environmental Protection Agency

- J Analyte was detected and the concentration is estimated.
- JB Concentration is estimated due to presence of blank contamination.
- U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

Seminar Name 1994-140 994-842 994-843 994-840 944-9 944-			Location Name		SP	2-B		SS-2B		SB-40			SB-	-41		SB-	-42
Margin CAMBO Consequent Campaign C				SP2-B-15.0	SP2-B-30.0	SP2-B-35.0	SP2-B-45.0		SB-40-29.5-30		SB-40-39-39.5	SB-41-24-25	SB-41-25.5-26	SB-41-28.5-29	SB-41-49-50	SB-42-25.5-26	SB-42-29-30
Mary Mary			Sample Date	6/11/2010	6/11/2010	6/11/2010	6/11/2010	6/22/2010	10/1/2018	10/1/2018	10/1/2018	10/2/2018	10/2/2018	10/2/2018	10/2/2018	10/3/2018	10/3/2018
Marcon			Sample Depth (feet)	15.0-15.0	30.0-30.0	35.0-35.0	45.0-45.0	1.0-1.0	29.5-30.0	31.5-32.0	39.0-39.5	24.0-25.0	25.5-26.0	28.5-29.0	49.0-50.0	25.5-26.0	29.0-30.0
According	Analyte	CAS No.	Screening Level (1)														1
Service	Metals																
Excellen	Antimony	7440-36-0	32														
Carbinary Carb	Arsenic	7440-38-2	20														
Comparison T-460-84 3,000 1	Beryllium	7440-41-7	160														1
Control Cont	Cadmium	7440-43-9	2.0														'
East	Chromium	7440-47-3	2,000					7.02									
More 749-97-49 7.00	Copper	7440-50-8	3,200														
Sicility	Lead	7439-92-1	250														
Selection	Mercury	7439-97-6	2.0														
Selection 772-3-92 400	Nickel	7440-02-0	1,600														
The Table	Selenium	7782-49-2	400														
Trace	Silver	7440-22-4	400														
Trace																	1
Total Petrolum Hydrocathons	Zinc	7440-66-6	24,000														1
Display Disp	Total Petroleum Hydrocarbons		<u> </u>														
Semioalistic Organic Compounds	·	DRO					l	I	1	1		I		1			
Sembolatic Organic Compounds Sembolation	Gasoline	GRO															
3.4 As friction period 3.2									<u> </u>					<u> </u>			
2.4.5 frictioloophenol 95.95.4 8,000		90-12-0						I									
2.4-Directophenol 120-83-2 240			8.000														
2.4-Dintrophenol 10:83-2 240																	_
2.4-Dintrophenol 105:67:9 1,600																	
2.4-Ditrophenol 51:28:5 160																	
2-Methylphenol 95-57-8																	_
2-Methylpaphthalene																	
2-Mitrophenol 95-48-7 4,000																	
2-Nitrophenol 88-75-5																	
4.6-Dinitro-o-cresol 534-52-1 6.4																	+
4-Chloro-3-methylphenol 59-50-7 8,000			6.4														+
4-Methylphenol 106-44-5 8,000																	
A-Nitrophenol 100-02-7																	
Acenaphthene 83-32-9 4,800 0.20 U 69 37 0.092 Acenaphthylene 208-96-8 0.63 4.2 2.0 U 0.010 U			1														
Acenaphthylene 208-96-8 0.63 4.2 2.0 U 0.010 U 0.010 U 0.010 U 0.015 U 0.010 U 0.015 U 0.045 U 1.4 10 0.15 0.046 U 1.1 0.06 U 0.05 U 0.046 U 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U 0.046 U 0.045 U 0.045 U 0.045 U 0.046 U 0.045 U 0.045 U 0.046 U 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U 0.046 U 0.045 U 0.045 U 0.045 U 0.045 U 0.045 U 0.045				0.20 11	69	37	0.092										
Anthracene 120-12-7 24,000 1.3 28 15 0.031																	
Benzo(a)anthracene 56-55-3	· '													1			
Benzo(a)pyrene 50-32-8 0.10 26 6.3 3.9 0.010 U 0.99 1.7 0.045 U 5.8 4.7 0.045 U 0.046 U 5.3 0.00 Benzo(b)fluoranthene 205-99-2 31 8.2 5.3 0.010 U 0.78 1.5 0.045 U 4.7 4.4 0.045 U 0.046 U 4.6 0.00 Benzo(k)fluoranthene 207-08-9 5.6 4.0 U 2.0 U 0.010 U 0.86 1.4 0.045 U 4.7 3.7 0.045 U 4.2 0.00 Chrysene 218-01-9 15 13 8.2 0.015 2.1 3.1 0.045 U 4.7 3.7 0.045 U 4.2 0.00 Dibenzo(a,h)anthracene 53-70-3 4.7 4.0 U 2.0 U 0.010 U 0.14 0.27 0.045 U 1.1 0.64 0.045 U 0.046 U 0.088 0.00 Indeno(1,2,3-c,d)pyrene 193-39-5 14 4.0 U </td <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td> </td> <td>2 1</td> <td>3.0</td> <td>0.045.11</td> <td>1/1</td> <td>10</td> <td>0.15</td> <td>0.046.11</td> <td>11</td> <td>0.046 U</td>								 	2 1	3.0	0.045.11	1/1	10	0.15	0.046.11	11	0.046 U
Benzo(b)fluoranthene 205-99-2 31 8.2 5.3 0.010 U 0.78 1.5 0.045 U 4.7 4.4 0.045 U 0.046 U 4.6 0.00 Benzo(k)fluoranthene 207-08-9 5.6 4.0 U 2.0 U 0.010 U 0.86 1.4 0.045 U 4.7 3.7 0.045 U 0.046 U 4.2 0.00 Chrysene 218-01-9 15 13 8.2 0.015 2.1 3.1 0.045 U 9.1 0.11 0.046 U 4.2 0.00 Dibenzo(a,h)anthracene 53-70-3 4.7 4.0 U 2.0 U 0.010 U 0.14 0.27 0.045 U 1.1 0.64 0.045 U 0.046 U 0.88 0.00 Indeno(1,2,3-c,d)pyrene 193-39-5 14 4.0 U 2.0 U 0.010 U 0.29 0.50 0.045 U 2.0 0.045 U 0.045 U 0.045 U 0.045 U 1.6 0.04 cPAHs (MTCA TEQ-HalfND) BaPEq	, ,				_												0.046 U
Benzo(k)fluoranthene 207-08-9 5.6 4.0 U 2.0 U 0.010 U 0.86 1.4 0.045 U 4.7 3.7 0.045 U 0.046 U 4.2 0.045 U Chrysene 218-01-9 15 13 8.2 0.015 2.1 3.1 0.045 U 12 9.1 0.11 0.046 U 11 0.04 Dibenzo(a,h)anthracene 53-70-3 4.7 4.0 U 2.0 U 0.010 U 0.14 0.27 0.045 U 1.1 0.046 U 0.046 U 0.88 0.04 Indeno(1,2,3-c,d)pyrene 193-39-5 14 4.0 U 2.0 U 0.010 U 0.29 0.50 0.045 U 2.0 0.046 U 1.6 0.04 CPAHs (MTCA TEQ-HalfND) BaPEq (U=1/2) 0.10 32 9.2 5.6 0.0086 1.4 2.4 0.045 U 8.6 6.8 0.046 U 7.7 0.04 CPAHs (MTCA TEQ-ZeroND) BaPEq (U=0) 0.10 32 8.6 5.3 <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td> </td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>+</td> <td></td> <td></td> <td>0.046 U</td>								 						+			0.046 U
Chrysene 218-01-9 15 13 8.2 0.015 2.1 3.1 0.045 U 12 9.1 0.11 0.046 U 11 0.046 U 11 0.046 U 12 0.045 U 12 0										<u> </u>							0.046 U
Dibenzo(a,h)anthracene 53-70-3 4.7 4.0 U 2.0 U 0.010 U 0.14 0.27 0.045 U 1.1 0.64 0.045 U 0.046 U 0.88 0.00 U Indeno(1,2,3-c,d)pyrene 193-39-5 14 4.0 U 2.0 U 0.010 U 0.29 0.50 0.045 U 2.0 1.2 0.045 U 0.046 U 1.6 0.04 cPAHs (MTCA TEQ-HalfND) BaPEq (U=1/2) 0.10 32 9.2 5.6 0.0086 1.4 2.4 0.045 U 8.6 6.8 0.046 U 7.7 0.04 cPAHs (MTCA TEQ-ZeroND) BaPEq (U=0) 0.10 32 8.6 5.3 0.0016 1.4 2.4 0.045 U 8.6 6.8 0.016 0.046 U 7.7 0.04 Benzo(g,h,i)perylene 191-24-2 13 4.0 U 2.0 U 0.010 U 0.010 U 0.045 U 8.6 6.8 0.016 U 7.7 0.04			+			1											0.046 U
Indeno(1,2,3-c,d)pyrene 193-39-5 14 4.0 U 2.0 U 0.010 U 0.29 0.50 0.045 U 2.0 1.2 0.045 U 0.046 U 1.6 0.046 U cPAHs (MTCA TEQ-HalfND) BaPEq (U=1/2) 0.10 32 9.2 5.6 0.0086 1.4 2.4 0.045 U 8.6 6.8 0.046 U 7.7 0.046 U cPAHs (MTCA TEQ-ZeroND) BaPEq (U=0) 0.10 32 8.6 5.3 0.0016 1.4 2.4 0.045 U 8.6 6.8 0.016 0.046 U 7.7 0.046 U Benzo(g,h,i)perylene 191-24-2 13 4.0 U 2.0 U 0.010 U 0.045 U 8.6 6.8 0.016 U 7.7 0.046 U																	
CPAHs (MTCA TEQ-HalfND) BaPEq (U=1/2) 0.10 32 9.2 5.6 0.0086 1.4 2.4 0.045 U 8.6 6.8 0.046 U 7.7 0.046 U cPAHs (MTCA TEQ-ZeroND) BaPEq (U=0) 0.10 32 8.6 5.3 0.0016 1.4 2.4 0.045 U 8.6 6.8 0.016 0.046 U 7.7 0.046 U Benzo(g,h,i)perylene 191-24-2 13 4.0 U 2.0 U 0.010 U 0.045 U 8.6 6.8 0.046 U 7.7 0.046 U														+			0.046 U
cPAHs (MTCA TEQ-ZeroND) BaPEq (U=0) 0.10 32 8.6 5.3 0.0016 1.4 2.4 0.045 U 8.6 6.8 0.016 0.046 U 7.7 0.045 U Benzo(g,h,i)perylene 191-24-2 13 4.0 U 2.0 U 0.010 U														+			0.046 U
Benzo(g,h,i)perylene 191-24-2 13 4.0 U 2.0 U 0.010 U																	0.046 U
						1			1.4	2.4	0.045 U	8.6	6.8	0.016	U.U46 U	1.1	0.046 U
Fluoranthene 206-44-0 3,200 0.94 54 32 0.057										1				 			



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name		SP2	2-B		SS-2B		SB-40			SB-	41		SB-	42
		Sample Name	SP2-B-15.0	SP2-B-30.0	SP2-B-35.0	SP2-B-45.0	SS-2B	SB-40-29.5-30	SB-40-31.5-32	SB-40-39-39.5	SB-41-24-25	SB-41-25.5-26	SB-41-28.5-29	SB-41-49-50	SB-42-25.5-26	SB-42-29-30
		Sample Date	6/11/2010	6/11/2010	6/11/2010	6/11/2010	6/22/2010	10/1/2018	10/1/2018	10/1/2018	10/2/2018	10/2/2018	10/2/2018	10/2/2018	10/3/2018	10/3/2018
		Sample Depth (feet)	15.0–15.0	30.0–30.0	35.0–35.0	45.0–45.0	1.0-1.0	29.5–30.0	31.5–32.0	39.0–39.5	24.0-25.0	25.5–26.0	28.5–29.0	49.0–50.0	25.5–26.0	29.0–30.0
Analyte	CAS No.	Screening Level (1)														
Semivolatile Organic Compounds	(cont.)															
Fluorene	86-73-7	3,200	0.20 U	61	34	0.055										
Naphthalene	91-20-3	5.0 (2)	0.20 U (2)	₄₇₀ (2)	160 (2)	0.13 (2)		88 1 (2)	140 J (2)	0.067 UJ (2)	160 J (2)	190 J (2)	0.063 UJ (2)	8.9 J (2)	130 J (2)	0.15 JB (2)
Pentachlorophenol	87-86-5	2.5	3.0 U	3.0 U	3.0 U	0.30 U										
Phenanthrene	85-01-8		0.49	140	82	0.14										
Phenol	108-95-2	24,000														
Pyrene	129-00-0	2,400	5.2	43	26	0.045										
Tetrachlorophenols (total)	25167-83-3															
Total Naphthalenes		5.0														
Volatile Organic Compounds	•															
1,1,1-Trichloroethane	71-55-6	2						0.037 UJ	0.034 UJ	0.034 UJ	0.022 UJ	0.024 UJ	0.032 UJ	0.028 UJ	0.028 UJ	0.031 UJ
1,1-Dichloroethane	75-34-3	180						0.03 UJ	0.027 UJ	0.027 UJ	0.041 UJ	0.045 UJ	0.025 UJ	0.022 UJ	0.054 UJ	0.025 UJ
1,2,4-Trimethylbenzene	95-63-6	1.3						2.4 J	4 J	0.027 UJ	2.3 J	4.2 J	0.025 UJ	0.050 J	2.1 J	0.025 UJ
1,2-Dichlorobenzene	95-50-1															
1,3,5-Trimethylbenzene	108-67-8	1.3						0.96 JB	2 J	0.034 UJ	0.076 UJ	0.083 UJ	0.032 UJ	0.028 UJ	1.1 J	0.031 UJ
1,3-Dichlorobenzene	541-73-1															
1,4-Dichlorobenzene	106-46-7															
Benzene	71-43-2	0.030						0.03 UJ	0.027 UJ	0.027 UJ	0.074 UJ	0.081 UJ	0.025 UJ	0.022 UJ	0.097 UJ	0.025 UJ
Chlorobenzene	108-90-7															
Chloroform	67-66-3	32						0.03 UJ	0.027 UJ	0.027 UJ	0.05 UJ	0.055 UJ	0.025 UJ	0.022 UJ	0.065 UJ	0.025 UJ
Chloromethane	74-87-3							0.074 UJ	0.067 UJ	0.067 UJ	0.027 UJ	0.03 UJ	0.063 UJ	0.056 UJ	0.036 UJ	0.062 UJ
cis-1,2-Dichloroethene	156-59-2	160						0.03 UJ	0.027 UJ	0.027 UJ	0.037 UJ	0.041 UJ	0.025 UJ	0.022 UJ	0.049 UJ	0.025 UJ
Ethylbenzene	100-41-4	6.0						0.20 J	0.14 J	0.034 UJ	0.10 UJ	0.25 J	0.032 UJ	0.11 J	0.13 UJ	0.031 UJ
Isopropylbenzene	98-82-8	8,000						0.18 J	0.36 J	0.034 UJ	0.13 J	0.32 J	0.032 UJ	0.031 J	0.13 UJ	0.031 UJ
Methyl isobutyl ketone	108-10-1	6,400						0.37 UJ	0.34 UJ	0.34 UJ	0.35 UJ	0.38 UJ	0.32 UJ	0.28 UJ	0.46 UJ	0.31 UJ
n-Propylbenzene	103-65-1	8,000						0.15 J	0.35 J	0.034 UJ	0.094 UJ	0.10 UJ	0.032 UJ	0.035 J	0.12 UJ	0.031 UJ
Styrene	100-42-5	16,000						0.13 J	0.064 J	0.034 UJ	0.044 UJ	0.048 UJ	0.032 UJ	0.028 UJ	0.058 UJ	0.031 UJ
Toluene	108-88-3	7.0						0.03 UJ	0.027 UJ	0.027 UJ	0.023 UJ	0.025 UJ	0.028 J	0.026 J	0.21 J	0.025 UJ
Trichloroethene	79-01-6	0.030						0.03 UJ	0.027 UJ	0.027 UJ	0.023 UJ	0.025 UJ	0.025 UJ	0.022 UJ	0.03 UJ	0.025 UJ
Xylene (meta & para)	108-38-3/106-42-3							0.80 J	0.52 J	0.067 UJ	0.15 J	0.64 J	0.063 UJ	0.056 UJ	0.059 UJ	0.062 UJ
Xylene (ortho)	95-47-6							0.32 J	0.20 J	0.034 UJ	0.079 UJ	0.35 J	0.032 UJ	0.041 J	0.10 UJ	0.031 UJ
Xylene (total)	1330-20-7	9.0						1.1 J	0.71 J	0.067 UJ	0.15 J	0.99 J	0.063 UJ	0.041 J	0.10 UJ	0.062 UJ

Blank cells are intentional.

All criteria and results are rounded to two significant figures.

-- Not available or not established.

BOLD Detected result exceeds criteria. The screening level is a proposed Site cleanup level.

- 1 Soil screening levels for all compounds designated as chemicals of interest are presented in Table 5.5. Landau Associates made no modifications to the analytical tables taken from February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.
- 2 Where naphthalene was run by USEPA 8260 only, or methylnaphthalenes were not run by USEPA 8270, the naphthalene result is compared to the screening level for total naphthalenes.
- 3 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

mg/kg = Milligrams per kilogram

MTCA = Model Toxics Control Act TEQ Toxic equivalent

USEPA = U.S. Environmental Protection Agency

- J Analyte was detected and the concentration is estimated.
- JB Concentration is estimated due to presence of blank contamination. U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name				SB-43					MV	V-14	
		Sample Name	SB-43-28.5-29	SB-43-31-31.5	SB-43-37.5-38	SB-143-37.5-38	SB-43-41-41.5	SB-43-47-47.5	SB-43-71-71.5	MW-14-20.5-21	MW-14-22-22.5	MW-14-27-28	MW-14-31-31
		Sample Date	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/5/2018	10/5/2018	10/5/2018	10/5/2018
		Sample Depth (feet)	28.5-29.0	31.0-31.5	37.5-38.0	37.5-38.0	41.0-41.5	47.0-47.5	71.0-71.5	20.5-21.0	22.0-22.5	27.0-28.0	31.0-31.5
Analyte	CAS No.	Screening Level (1)											
Metals													
Antimony	7440-36-0	32		0.18 U									
Arsenic	7440-38-2	20		1.4									
Beryllium	7440-41-7	160		0.18 U									
Cadmium	7440-43-9	2.0		0.18 U									
Chromium	7440-47-3	2,000		23									
Copper	7440-50-8	3,200		9.9									
Lead	7439-92-1	250		1.4									
Mercury	7439-97-6	2.0		0.29 U									
Nickel	7440-02-0	1,600		34									
Selenium	7782-49-2	400		0.83									
Silver	7440-22-4	400		0.092 U									
Thallium	7440-28-0	0.80		0.18 U									
Zinc	7440-66-6	24,000		25									
otal Petroleum Hydrocarbons		,											
Diesel	DRO				1								1
Gasoline	GRO												
emivolatile Organic Compounds													
1-Methylnaphthalene	90-12-0			160									
2,4,5-Trichlorophenol	95-95-4	8,000		0.11 U									
2,4,6-Trichlorophenol	88-06-2	80		0.11 U									
2,4-Dichlorophenol	120-83-2	240		0.11 U									
2,4-Dimethylphenol	105-67-9	1,600		0.19									_
2,4-Dinitrophenol	51-28-5	160		0.56 UJ									
2-Chlorophenol	95-57-8	400		0.11 U									+
2-Methylnaphthalene	91-57-6			290									+
2-Methylphenol	95-48-7	4,000		0.12									
2-Nitrophenol	88-75-5	4,000		0.12 0.11 U									
4,6-Dinitro-o-cresol	534-52-1	6.4		0.21 U									
4-Chloro-3-methylphenol	59-50-7	8,000		0.21 U									
4-Methylphenol	106-44-5	8,000		0.21 U							 		+
4-Nitrophenol	100-02-7			67	1						 		+
Acenaphthene	83-32-9	4,800		140	1						 		+
· · · · · · · · · · · · · · · · · · ·	208-96-8	4,800		7.7							 		
Acenaphthylene Anthracene	208-96-8 120-12-7	24,000		51							-		+
Benzo(a)anthracene	56-55-3	24,000	5.5	26	2.4	2.2	0.040 U	0.048 U	0.047 U	14 J	0.046 U	0.060	0.040 U
, ,													
Benzo(a)pyrene	50-32-8	0.10	2.8	12	0.97	1.1	0.040 U	0.048 U	0.047 U	5.9 J	0.046 U	0.043 U	0.040 U
Benzo(b)fluoranthene	205-99-2		2.2	11	0.92	0.97	0.040 U	0.048 U	0.047 U	4.3 J	0.046 U	0.043 U	0.040 U
Benzo(k)fluoranthene	207-08-9		2.5	9.7	0.70	0.67	0.040 U	0.048 U	0.047 U	4.0 J	0.046 U	0.043 U	0.040 U
Chrysene	218-01-9		5.5	25	2.3	2.2	0.040 U	0.048 U	0.047 U	14 J	0.046 U	0.062	0.040 U
Dibenzo(a,h)anthracene	53-70-3		0.46	2.4	0.21	0.2	0.040 U	0.048 U	0.047 U	0.71 J	0.046 U	0.043 U	0.040 U
Indeno(1,2,3-c,d)pyrene	193-39-5		0.84	3.2	0.40	0.35	0.040 U	0.048 U	0.047 U	1.5 J	0.046 U	0.043 U	0.040 U
cPAHs (MTCA TEQ-HalfND)	BaPEq (U=1/2)	0.10	4.0	17	1.5	1.5	0.040 U	0.048 U	0.047 U	8.4 J	0.046 U	0.036	0.040 U
cPAHs (MTCA TEQ-ZeroND)	BaPEq (U=0)	0.10	4.0	17	1.5	1.5	0.040 U	0.048 U	0.047 U	8.4 J	0.046 U	0.0066	0.040 U
Benzo(g,h,i)perylene	191-24-2			3.4									
Fluoranthene	206-44-0	3,200		110	ĺ		Ī		Ī]			1



Table 5.3 Soil Analytical Data 1992–2018 (mg/kg)

		Location Name				SB-43					MV	V-14	
		Sample Name	SR-43-28 5-29	SB-43-31-31.5	SB-43-37.5-38	SB-143-37.5-38	SB-43-41-41.5	SB-43-47-47.5	SR-43-71-71 5	MW-14-20.5-21	MW-14-22-22.5		MW-14-31-31.5
		Sample Date	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/4/2018	10/5/2018	10/5/2018	10/5/2018	10/5/2018
		Sample Depth (feet)	28.5–29.0	31.0-31.5	37.5–38.0	37.5–38.0	41.0-41.5	47.0-47.5	71.0-71.5	20.5–21.0	22.0–22.5	27.0–28.0	31.0-31.5
Analyte	CAS No.	Screening Level (1)	20.3 23.0	31.0 31.3	37.3 30.0	37.3 30.0	41.0 41.5	47.0 47.5	71.0 71.5	20.5 21.0	22.0 22.3	27.0 20.0	31.0 31.5
Semivolatile Organic Compounds													
Fluorene	86-73-7	3,200		110	T		I	I				l	T
Naphthalene	91-20-3	5.0 ⁽²⁾	540 ⁽²⁾	890 ⁽³⁾	78 ⁽²⁾	52 ⁽²⁾	0.048 UJ ⁽²⁾	13 ⁽²⁾	0.059 UJ ⁽²⁾	170 ⁽²⁾	0.19 (2)	0.54 ⁽²⁾	1.8 (2)
Pentachlorophenol	87-86-5	2.5	0.0	0.23 J	1	<u> </u>	0.0.0		0.000		0.23	0.5 .	1 2.0
Phenanthrene	85-01-8			310									•
Phenol	108-95-2	24,000		0.11 U									•
Pyrene	129-00-0	2,400		97									
Tetrachlorophenols (total)	25167-83-3												
Total Naphthalenes		5.0		1,300									
Volatile Organic Compounds	<u>. </u>				•		•	•					
1,1,1-Trichloroethane	71-55-6	2	0.024 U	0.024 U	0.022 U	0.027 U	0.024 U	0.03 U	0.029 U	0.021 U	0.026 U	0.027 U	0.036 U
1,1-Dichloroethane	75-34-3	180	0.045 U	0.046 U	0.041 U	0.021 U	0.019 U	0.024 U	0.024 U	0.041 U	0.021 U	0.022 U	0.029 U
1,2,4-Trimethylbenzene	95-63-6	1.3	13	34	1.9	1.5	0.019 U	0.30	0.024 U	0.43	0.021 U	0.022 U	0.029 U
1,2-Dichlorobenzene	95-50-1												
1,3,5-Trimethylbenzene	108-67-8	1.3	0.083 U	13	0.076 U	0.57	0.024 U	0.11	0.029 U	0.075 U	0.026 U	0.027 U	0.036 U
1,3-Dichlorobenzene	541-73-1												
1,4-Dichlorobenzene	106-46-7												
Benzene	71-43-2	0.030	0.081 U	0.083 U	0.074 U	0.021 U	0.019 U	0.024 U	0.024 U	0.073 U	0.021 U	0.022 U	0.029 U
Chlorobenzene	108-90-7												
Chloroform	67-66-3	32	0.054 U	0.056 U	0.050 U	0.021 U	0.019 U	0.024 U	0.024 U	0.049 U	0.021 U	0.022 U	0.029 U
Chloromethane	74-87-3		0.030 U	0.031 U	0.027 U	0.053 U	0.048 U	0.059 U	0.059 U	0.027 U	0.052 U	0.054 U	0.071 U
cis-1,2-Dichloroethene	156-59-2	160	0.040 U	0.042 U	0.037 U	0.021 U	0.019 U	0.024 U	0.024 U	0.037 U	0.021 U	0.022 U	0.029 U
Ethylbenzene	100-41-4	6.0	2.9	13	0.53	0.44	0.024 U	0.46	0.029 U	0.10 U	0.026 U	0.027 U	0.036 U
Isopropylbenzene	98-82-8	8,000	1.2	3.3	0.19 J	0.13	0.024 U	0.059	0.029 U	0.096 U	0.026 U	0.027 U	0.036 U
Methyl isobutyl ketone	108-10-1	6,400	0.38 U	0.39 U	0.35 U	0.27 U	0.24 U	0.30 U	0.29 U	0.35 U	0.26 U	0.27 U	0.36 U
n-Propylbenzene	103-65-1	8,000	0.91	2.5	0.094 U	0.097	0.024 U	0.058	0.029 U	0.093 U	0.026 U	0.027 U	0.036 U
Styrene	100-42-5	16,000	2.4	11	0.12 J	0.027 U	0.024 U	0.030 U	0.029 U	0.043 U	0.026 U	0.027 U	0.036 U
Toluene	108-88-3	7.0	0.23 J	0.65	0.023 U	0.021 U	0.019 U	0.035	0.024 U	0.14 J	0.021 U	0.022 U	0.029 U
Trichloroethene	79-01-6	0.030	0.025 U	0.026 U	0.023 U	0.021 U	0.019 U	0.024 U	0.024 U	0.023 U	0.021 U	0.022 U	0.029 U
Xylene (meta & para)	108-38-3/106-42-3		13	42	0.89	0.72	0.048 U	0.16	0.059 U	0.044 U	0.052 U	0.054 U	0.071 U
Xylene (ortho)	95-47-6		5.1	18	0.45	0.35	0.024 U	0.14	0.029 U	0.078 U	0.026 U	0.027 U	0.036 U
Xylene (total)	1330-20-7	9.0	18	60	1.3	1.1	0.048 U	0.30	0.059 U	0.078 U	0.052 U	0.054 U	0.071 U

Blank cells are intentional.

All criteria and results are rounded to two significant figures.

-- Not available or not established.

BOLD Detected result exceeds criteria. The screening level is a proposed Site cleanup level.

- 1 Soil screening levels for all compounds designated as chemicals of interest are presented in Table 5.5. Landau Associates made no modifications to the analytical tables taken from the February 2022 Remedial Investigation and Feasibility Study apart from updating screening levels.
- 2 Where naphthalene was run by USEPA 8260 only, or methylnaphthalenes were not run by USEPA 8270, the naphthalene result is compared to the screening level for total naphthalenes.
- 3 Naphthalene was analyzed by both USEPA 8260 and 8270. The 8270 result was preferred and is shown here and the calculated total naphthalenes concentration is compared to the screening level.

Abbreviations:

CAS = Chemical Abstracts Service

cPAH = Carcinogenic polycyclic aromatic hydrocarbon

mg/kg = Milligrams per kilogram

MTCA = Model Toxics Control Act TEQ Toxic equivalent

USEPA = U.S. Environmental Protection Agency

- J Analyte was detected and the concentration is estimated.
- JB Concentration is estimated due to presence of blank contamination. U Analyte is not detected at the associated reporting limit.
- UJ Analyte is not detected at the associated reporting limit, which is an estimate.